

## EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	86	(torre buezo jadhav mitch pedregal tercero).in. and ("546" or "548").clas.	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/10/17 14:30
L3	8	I1 and opioid	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/10/17 14:30

=> b reg  
FILE 'REGISTRY' ENTERED AT 13:16:50 ON 17 OCT 2007  
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STRUCTURE FILE UPDATES: 16 OCT 2007 HIGHEST RN 950817-67-1  
DICTIONARY FILE UPDATES: 16 OCT 2007 HIGHEST RN 950817-67-1

New CAS Information Use Policies, enter HELP USAGETERMS for details.

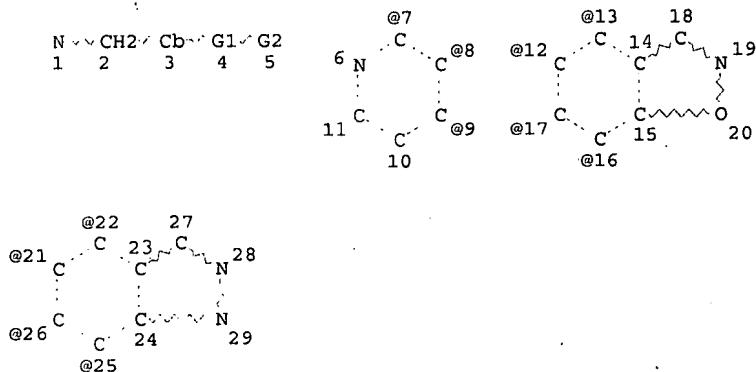
TSCA INFORMATION NOW CURRENT THROUGH June 29, 2007

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conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and  
predicted properties as well as tags indicating availability of  
experimental property data in the original document. For information  
on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=> d que sta 122  
L16 2025948 SEA FILE=REGISTRY ABB=ON PLU=ON 46.150.18/RID AND (NC5 OR  
NOC3-C6 OR N2C3-C6)/ES  
L20 STR



VAR G1=O/S/SE  
VAR G2=7/8/9/12/13/16/17/22/21/25/26

NODE ATTRIBUTES:

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GGCAT IS MCY UNS AT 3  
DEFAULT ECLEVEL IS LIMITED  
ECOUNT IS E6 C AT 3

GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED  
NUMBER OF NODES IS 29

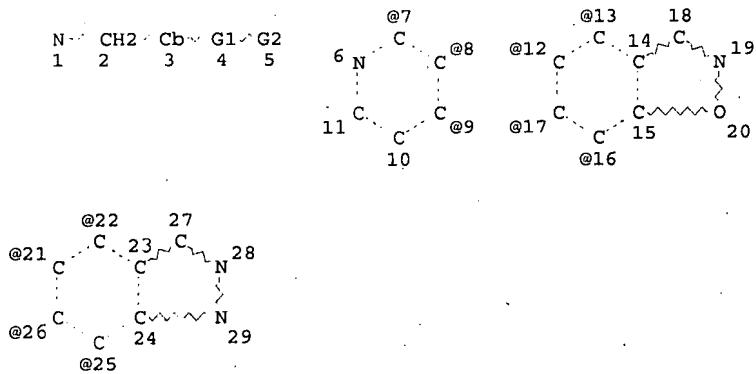
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L22 679 SEA FILE=REGISTRY SUB=L16 SSS FUL L20

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SEARCH TIME: 00.00.07

679 ANSWERS

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NOC3-C6 OR N2C3-C6)/ES  
L20 STR



VAR G1=O/S/SE

VAR G2=7/8/9/12/13/16/17/22/21/25/26

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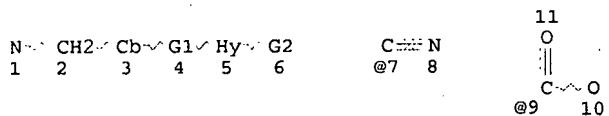
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STEREO ATTRIBUTES: NONE

L22 679 SEA FILE=REGISTRY SUB=L16 SSS FUL L20  
L35 STR

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VAR G2=7/9

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DEFAULT ECLEVEL IS LIMITED

ECOUNT IS E6 C AT 3

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STEREO ATTRIBUTES: NONE

L37 276 SEA FILE=REGISTRY SUB=L22 SSS FUL L35

100.0% PROCESSED 679 ITERATIONS  
SEARCH TIME: 00.00.01

276 ANSWERS

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 FILE 'HCAPLUS' ENTERED AT 13:17:26 ON 17 OCT 2007  
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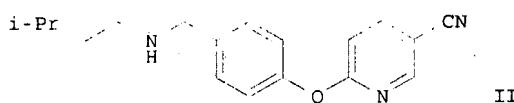
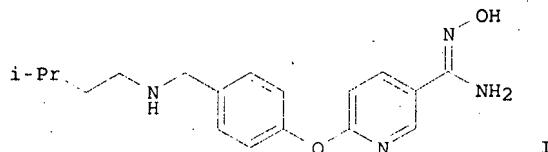
FILE COVERS 1907 - 17 Oct 2007 VOL 147 ISS 17  
FILE LAST UPDATED: 16 Oct 2007 (20071016/ED)

New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> => d bib abs hitstr 146

L46 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2007 ACS on STN  
AN 2005:638874 HCAPLUS  
DN 143:153386  
TI A preparation of di(hetero)aryl ethers, useful as opioid receptor antagonists  
IN De La Torre, Marta Garcia; Diaz Buezo, Nuria; Jadhav, Prabhakar Kondaji; Mitch, Charles Howard; Pedregal-Tercero, Concepcion  
PA Eli Lilly and Company, USA  
SO PCT Int. Appl., 67 pp.  
CODEN: PIXXD2  
DT Patent  
LA English  
FAN.CNT 1  
PATENT NO. KIND DATE APPLICATION NO. DATE  
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PI WO2005066164 A1 20050721 2004WO-US39766 20041215  
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GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC,  
LK, LR, LS, LT, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI,  
NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY,  
TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW  
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MR, NE, SN, TD, TG  
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CA--2549089 A1 20050721 2004CA-2549089 20041215  
EP--1699783 A1 20060913 2004EP-0812312 20041215  
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BR2004017714 A 20070320 2004BR-0017714 20041215  
JP2007516284 T 20070621 2006JP-0547016 20041215  
US2007112036 A1 20070517 2006US-0581178 20060531  
IN2006DN03118 A 20070824 2006IN-DN03118 20060531  
MX2006PA07198 A 20060904 2006MX-PA07198 20060622  
PRAI 2003EP-0380303 A 20031222  
2004US-539748P P 20040128  
2004WO-US39766 W 20041215  
OS CASREACT 143:153386; MARPAT 143:153386  
GI



AB The invention relates to a preparation of di(hetero)aryl ethers, useful as opioid receptor antagonists. The invention compds. are useful for the

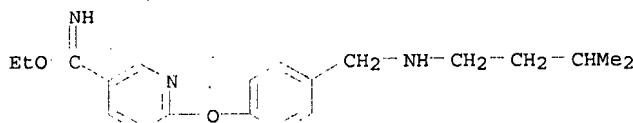
treatment, prevention or amelioration of obesity and related diseases. For instance, phenoxy nicotinamide derivative I [GTP- $\gamma$ -S binding assay,  $K_i$  (nM):  $\mu$  - 43.43,  $\kappa$  - 117.09,  $\delta$  - 269.06] was prepared via addition of hydroxylamine to phenoxy nicotinonitrile derivative II with a yield of 63%.

IT 859505-27-4P

RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)  
(preparation of di(hetero)aryl ethers useful as opioid receptor antagonists)

RN 859505-27-4 HCPLUS

CN 3-Pyridinecarboximidic acid, 6-[4-[[[(3-methylbutyl)amino]methyl]phenoxy]-, ethyl ester (CA INDEX NAME)



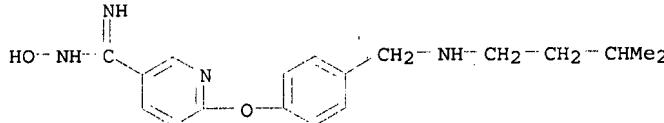
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859505-31-0P 859505-33-2P 859505-34-3P  
859505-35-4P 859505-36-5P 859505-37-6P  
859505-39-8P 859505-47-8P 859505-48-9P  
859505-49-0P 859505-50-3P 859505-52-5P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(preparation of di(hetero)aryl ethers useful as opioid receptor antagonists)

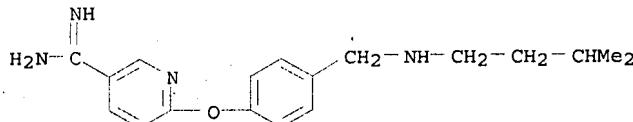
RN 859505-28-5 HCPLUS

CN 3-Pyridinecarboximidamide, N-hydroxy-6-[4-[[[(3-methylbutyl)amino]methyl]phenoxy]- (CA INDEX NAME)



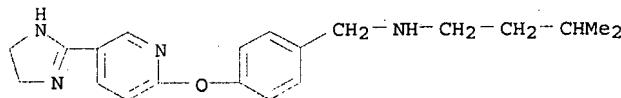
RN 859505-29-6 HCPLUS

CN 3-Pyridinecarboximidamide, 6-[4-[[[(3-methylbutyl)amino]methyl]phenoxy]- (CA INDEX NAME)



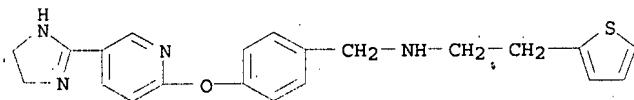
RN 859505-30-9 HCPLUS

CN Benzenemethanamine, 4-[[5-(4,5-dihydro-1H-imidazol-2-yl)-2-pyridinyl]oxy]-N-(3-methylbutyl)- (CA INDEX NAME)

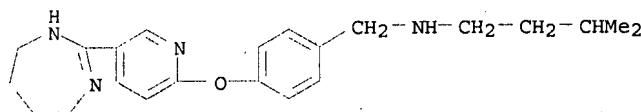


RN 859505-31-0 HCPLUS

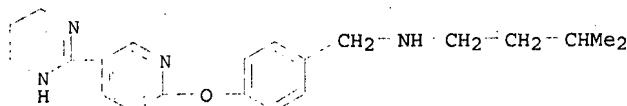
CN 2-Thiopheneethanamine, N-[[4-[[5-(4,5-dihydro-1H-imidazol-2-yl)-2-pyridinyl]oxy]phenyl]methyl]- (CA INDEX NAME)



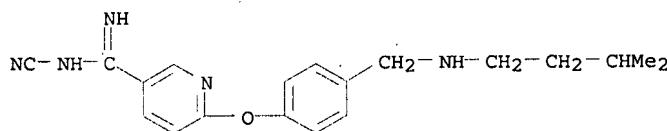
RN 859505-33-2 HCPLUS  
 CN Benzenemethanamine, N-(3-methylbutyl)-4-[[5-(4,5,6,7-tetrahydro-1H-1,3-diazepin-2-yl)-2-pyridinyl]oxy]- (CA INDEX NAME)



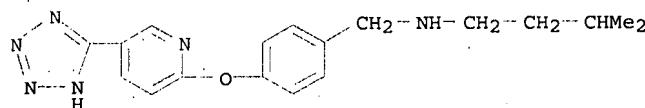
RN 859505-34-3 HCPLUS  
 CN Benzenemethanamine, N-(3-methylbutyl)-4-[[5-(1,4,5,6-tetrahydro-2-pyrimidinyl)-2-pyridinyl]oxy]- (CA INDEX NAME)



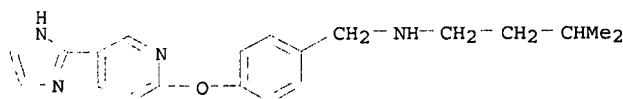
RN 859505-35-4 HCPLUS  
 CN 3-Pyridinecarboximidamide, N-cyano-6-[4-[(3-methylbutyl)amino]methyl]phenoxy]- (CA INDEX NAME)



RN 859505-36-5 HCPLUS  
 CN Benzenemethanamine, N-(3-methylbutyl)-4-[[5-(1H-tetrazol-5-yl)-2-pyridinyl]oxy]- (9CI) (CA INDEX NAME)

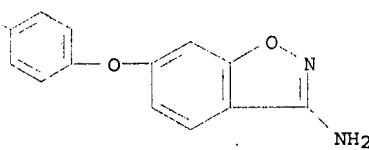


RN 859505-37-6 HCPLUS  
 CN Benzenemethanamine, 4-[[5-(1H-imidazol-2-yl)-2-pyridinyl]oxy]-N-(3-methylbutyl)- (CA INDEX NAME)



RN 859505-39-8 HCPLUS  
 CN 1,2-Benzisoxazol-3-amine, 6-[[4-[(3-methylbutyl)amino]methyl]phenoxy]- (CA INDEX NAME)

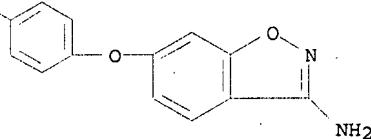
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RN 859505-47-8 HCPLUS

CN 1,2-Benzisoxazol-3-amine, 6-[4-[(3,3-dimethylbutyl)amino]methyl]phenoxy]-  
(CA INDEX NAME)

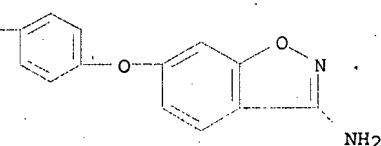
Me<sub>3</sub>C—CH<sub>2</sub>—CH<sub>2</sub>—NH—CH<sub>2</sub>



RN 859505-48-9 HCPLUS

CN 1,2-Benzisoxazol-3-amine, 6-[4-[(2-(4-fluorophenyl)ethyl)amino]methyl]phenoxy]-  
(CA INDEX NAME)

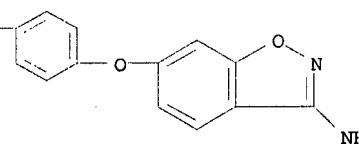
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RN 859505-49-0 HCPLUS

CN 1,2-Benzisoxazol-3-amine, 6-[4-[(2-(tetrahydro-2H-pyran-4-yl)ethyl)amino]methyl]phenoxy]-  
(CA INDEX NAME)

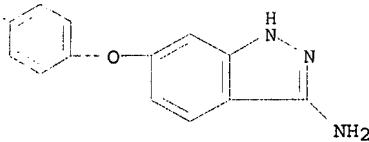
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RN 859505-50-3 HCPLUS

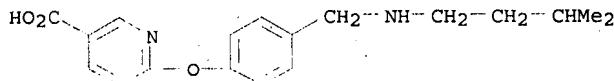
CN 1H-Indazol-3-amine, 6-[4-[(3-methylbutyl)amino]methyl]phenoxy]-  
(CA INDEX NAME)

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RN 859505-52-5 HCPLUS

CN 3-Pyridinecarboxylic acid, 6-[4-[(3-methylbutyl)amino]methyl]phenoxy]-  
(CA INDEX NAME)

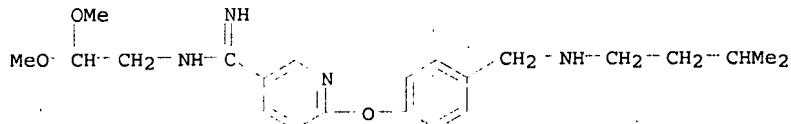


IT 859505-38-7P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
(preparation of di(hetero)aryl ethers useful as opioid receptor antagonists)

RN 859505-38-7 HCPLUS

CN 3-Pyridinecarboximidamide, N-(2,2-dimethoxyethyl)-6-[4-[(3-methylbutyl)amino]methyl]phenoxy - (CA INDEX NAME)

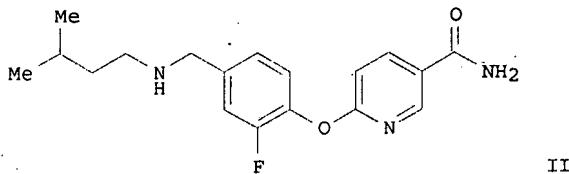
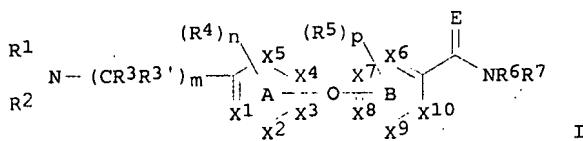


RE.CNT 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=&gt; d bib abs hitstr 145

L45 ANSWER 1 OF 1 HCPLUS COPYRIGHT 2007 ACS on STN  
AN 2004:267241 HCPLUS  
DN 140:303538  
TI Preparation of [(aminoalkyl)aryl]oxy]nicotinamides and analogs as opioid receptor antagonist for treatment of obesity and related conditions  
IN Blanco-Pillado, Maria-Jesus; Chappell, Mark Donald; Garcia De la Torre, Marta; Diaz Buezo, Nuria; Fritz, James Erwin; Holloway, William Glen; Matt, James Edward, Jr.; Mitch, Charles Howard; Pedregal-Tercero, Concepcion; Quimby, Steven James; Siegel, Miles Goodman; Smith, Dana Rae; Stucky, Russell Dean; Takeuchi, Kumiko; Thomas, Elizabeth Marie; Wolfe, Chad Nolan  
PA Eli Lilly and Company, USA  
SO PCT Int. Appl., 559 pp.  
CODEN: PIXXD2  
DT Patent  
LA English  
FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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PRAI 2002US-412158P	P	20020919 <--		
2003WO-US26300	W	20030917 <--		
OS MARPAT 140:303538				
GI				



AB Title diaryl ethers I [wherein X1-X10 = independently C, CH, or N; provided that each of rings A or B has no more than 2 N atoms; E = O or NH; R1 and R2 = independently H or (un)substituted (cyclo)alkyl, alkenyl, alkynyl, (alkyl)aryl, (aryl)heterocyclyl, (cyclo)alkylheterocyclyl, (cyclo)alkanoylalkyl, aroylalkyl, aryloxyalkyl, benzhydryl, bicycyl(alkyl), benzoyl(alkyl), alkoxyalkyl, alkoxy carbonyl, (aryl)alkylsulfonyl, heterocyclylalkylsulfonyl, cycloalkylalkyl, carboxyalkyl, carbamoylalkyl, etc.; R3 and R3' = independently H, alkyl, alkenyl, alkynyl, (alkyl)aryl, or alkylcycloalkyl; R4 and R5 = independently H, (halo)alkyl, alkenyl, alkynyl, alkoxy(halo)alkyl, thioalkyl, halo, aryl(alkyl), alkanoyl, alkoxy carbonyl, aminoalkyl, cycloalkylalkyl, etc.; R6 and R7 = independently H, (cyclo)alkyl, alkenyl, alkynyl, alkanoyl, OH, alkoxy, (aryl)alkylsulfonyl, heterocyclylalkylsulfonyl, aryl(alkyl), carbamoyl(alkyl), etc.; m = 1-3; n = 0-3; p = 0-3; or pharmaceutically acceptable salts, solvates, enantiomers, racemates, diastereomers, or mixts. thereof] were prepared as  $\mu$ -,  $\kappa$ -, and  $\delta$ -opioid receptor antagonists. For example, reductive amination of 6-(2-fluoro-4-formylphenoxy)nicotinamide and 3-methylbutylamine provided II (99%). The latter inhibited ex vivo binding of [<sup>3</sup>H]-diprenorphine in rat striatum/nucleus accumbens by >65% at a concentration of 7 mg/kg. In an acute feeding rat obesity assay, II suppressed opioid receptors at a dose of 0.3  $\mu$ g/kg. In addition, diet-induced obese rats achieved an energy balance (caloric intake minus utilization) of -81 kcal/kg/day upon administration of 0.3 mg/kg p.o. of II in an indirect calorimetry assay. Thus, I and their pharmaceutical compns. are useful for the treatment, prevention, or amelioration of obesity and related diseases.

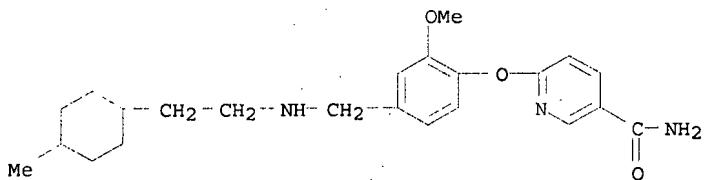
IT 676496-11-0P, 6-[2-Methoxy-4-[[[2-(4-methylcyclohexyl)ethyl]amino]methyl]phenoxy]nicotinamide 676496-29-0P, [4-[(6-Carbamoylpyridin-3-yl)oxy]benzyl](phenethyl)carbamic acid tert-butyl ester 676496-49-4P, 5-[4-[(3-Methylbutylamino)methyl]phenoxy]pyridine-2-carboxamide 676497-19-1P, 6-[2-Methoxy-4-[(4-methylpentylamino)methyl]phenoxy]nicotinamide 676497-21-5P, 6-[2-Methoxy-4-[[2-(p-tolyl)ethyl]amino]methyl]phenoxy]nicotinamide 676497-85-1P, 676501-30-7P, 5-[2-Methoxy-4-[[2-(tetrahydropyran-4-yl)ethyl]amino]methyl]phenoxy]pyridine-2-carboxamide 676501-43-2P, 6-[4-[(Phenethylamino)methyl]phenoxy]nicotinic acid ethyl ester 676501-44-3P, [4-[(5-Ethylcarbamoylpyridin-2-yl)oxy]benzyl](phenethyl)carbamic acid tert-butyl ester 676501-45-4P, 6-[4-[(tert-Butoxycarbonyl)(phenethyl)amino]methyl]phenoxy]nicotinic acid 676501-50-1P, 6-[4-[(3-Methylbutylamino)methyl]phenoxy]nicotinic acid ethyl ester 676501-51-2P, 6-[4-[(tert-Butoxycarbonyl)(phenethyl)amino]methyl]phenoxy]nicotinic acid ethyl ester 676501-52-3P, 6-[4-[(tert-Butoxycarbonyl)(3-methylbutyl)amino]methyl]phenoxy]nicotinic acid ethyl ester 676501-53-4P, 6-[4-[(tert-Butoxycarbonyl)(3-methylbutyl)amino]methyl]phenoxy]nicotinic acid 676501-54-5P, [4-[(5-(Isopropylcarbamoyl)pyridin-2-yl)oxy]benzyl](phenethyl)carbamic acid tert-butyl ester 676501-55-6P, N-(3-Methylbutyl)-N-[4-[(5-methylcarbamoylpyridin-2-yl)oxy]benzyl]carbamic acid tert-butyl ester 676501-56-7P, N-[4-[(5-Ethylcarbamoylpyridin-2-yl)oxy]benzyl](3-methylbutyl)carbamic acid tert-butyl ester 676501-57-8P, [4-[(5-(Isopropylcarbamoyl)pyridin-2-yl)oxy]benzyl](3-methylbutyl)carbamic acid tert-butyl ester

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(intermediate; preparation of (aryloxy)nicotinamides and analogs as opioid receptor antagonist for treatment of obesity and related conditions)

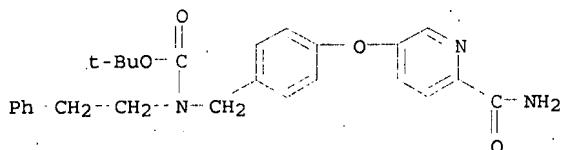
RN 676496-11-0 HCAPLUS

CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[[2-(4-methylcyclohexyl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



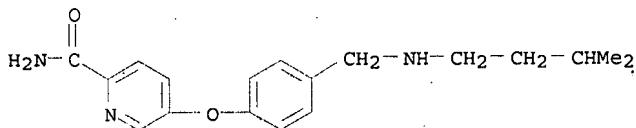
RN 676496-29-0 HCAPLUS

CN Carbamic acid, [[4-[[6-(aminocarbonyl)-3-pyridinyl]oxy]phenyl]methyl] (2-phenylethyl)-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)



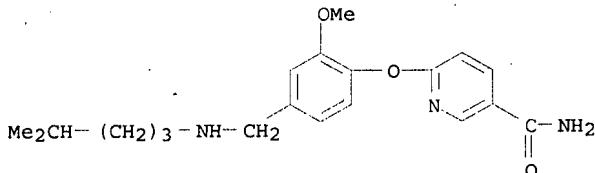
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CN 2-Pyridinecarboxamide, 5-[4-[[[(3-methylbutyl)amino]methyl]phenoxy] - (CA INDEX NAME)



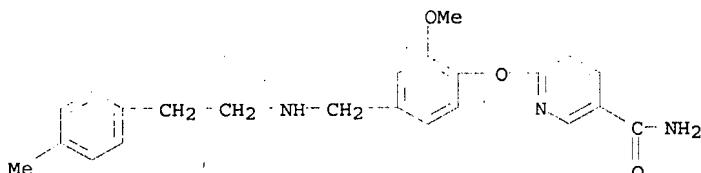
RN 676497-19-1 HCAPLUS

CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[[(4-methylpentyl)amino]methyl]phenoxy] - (CA INDEX NAME)



RN 676497-21-5 HCAPLUS

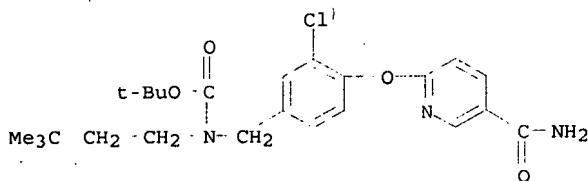
CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[[2-(4-methylphenyl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



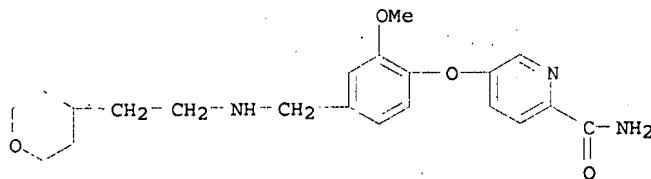
RN 676497-85-1 HCAPLUS

CN Carbamic acid, [[4-[[5-(aminocarbonyl)-2-pyridinyl]oxy] - 3 -

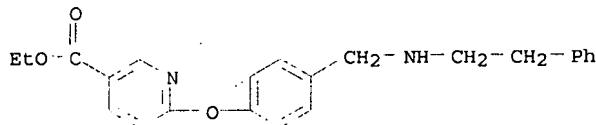
chlorophenyl]methyl] (3,3-dimethylbutyl)-, 1,1-dimethylethyl ester (9CI)  
(CA INDEX NAME)



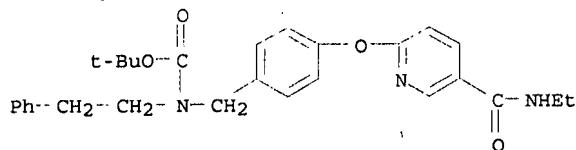
RN 676501-30-7 HCPLUS  
CN 2-Pyridinecarboxamide, 5-[2-methoxy-4-[[2-(tetrahydro-2H-pyran-4-yl)ethyl]amino]methyl]phenoxy]- (CA INDEX NAME)



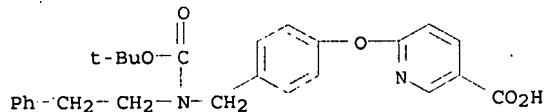
RN 676501-43-2 HCPLUS  
CN 3-Pyridinecarboxylic acid, 6-[4-[(2-phenylethyl)amino]methyl]phenoxy]-, ethyl ester (CA INDEX NAME)



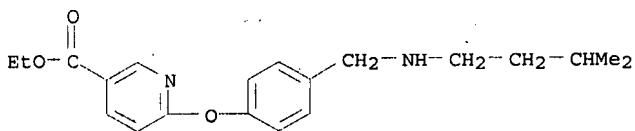
RN 676501-44-3 HCPLUS  
CN Carbamic acid, [[4-[[5-[(ethylamino)carbonyl]-2-pyridinyl]oxy]phenyl]methyl](2-phenylethyl)-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)



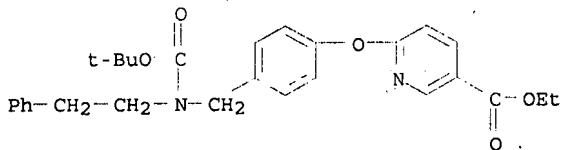
RN 676501-45-4 HCPLUS  
CN 3-Pyridinecarboxylic acid, 6-[4-[[[(1,1-dimethylethoxy)carbonyl](2-phenylethyl)amino]methyl]phenoxy]- (CA INDEX NAME)



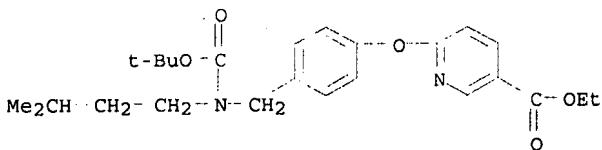
RN 676501-50-1 HCPLUS  
CN 3-Pyridinecarboxylic acid, 6-[4-[[[(3-methylbutyl)amino]methyl]phenoxy]-, ethyl ester (CA INDEX NAME)



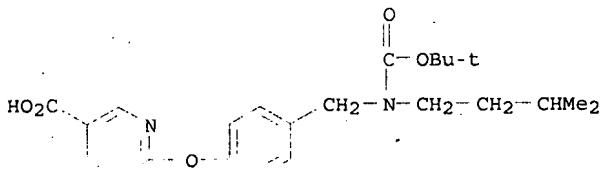
RN 676501-51-2 HCAPLUS  
 CN 3-Pyridinecarboxylic acid, 6-[4-[[[(1,1-dimethylethoxy)carbonyl](2-phenylethyl)amino]methyl]phenoxy]-, ethyl ester (CA INDEX NAME)



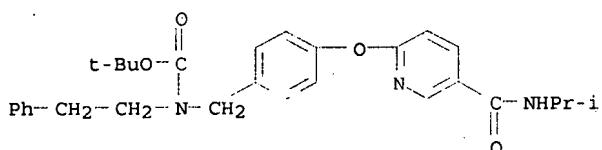
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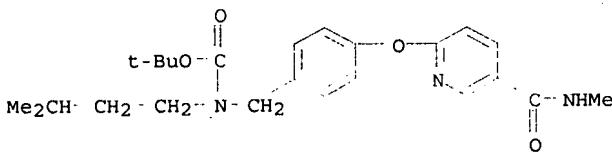
RN 676501-53-4 HCAPLUS  
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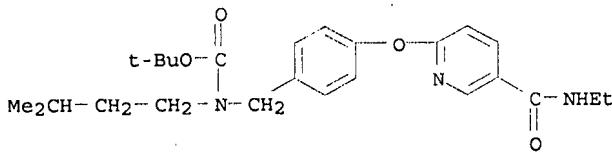
RN 676501-54-5 HCAPLUS  
 CN Carbamic acid, [[4-[[5-[(1-methylethyl)amino]carbonyl]-2-pyridinyl]oxy]phenyl]methyl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)



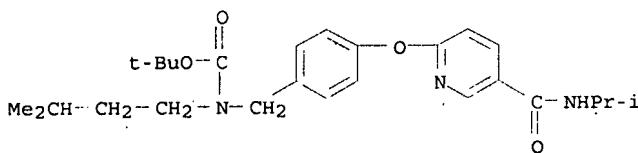
RN 676501-55-6 HCAPLUS  
 CN Carbamic acid, [[4-[[5-[(methylamino)carbonyl]-2-pyridinyl]oxy]phenyl]methyl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)



RN 676501-56-7 HCPLUS  
 CN Carbamic acid, [(4-[(5-[(ethylamino)carbonyl]-2-pyridinyl]oxy)phenyl]methyl](3-methylbutyl)-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

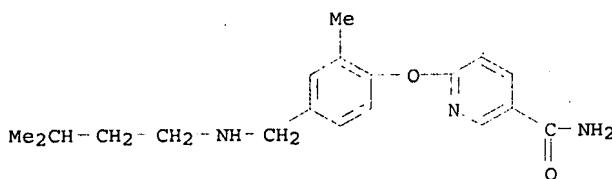


RN 676501-57-8 HCPLUS  
 CN Carbamic acid, (3-methylbutyl)[[4-[(5-[(1-methylethyl)amino]carbonyl]-2-pyridinyl]oxy]phenyl]methyl ester (9CI) (CA INDEX NAME)

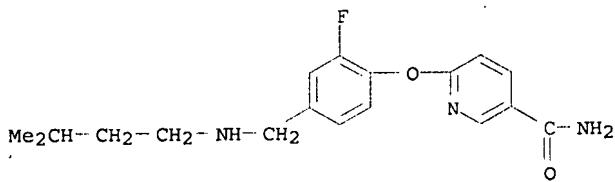


IT 676494-38-5P, 6-[2-Methyl-4-[(3-methylbutylamino)methyl]phenoxy]nicotinamide 676494-39-6P, 6-[2-Fluoro-4-[(3-methylbutyl)amino]methyl]phenoxy]nicotinamide 676497-13-5P, 6-[2-Methoxy-4-[(2-(tetrahydropyran-4-yl)ethyl)amino]methyl]phenoxy]nicotinamide 676497-16-8P, 6-[4-[(2-(4-Fluorophenyl)ethyl)amino]methyl-2-methoxyphenoxy]nicotinamide 676497-17-9P, 6-[4-[(2-(2-Fluorophenyl)ethyl)amino]methyl]-2-methoxyphenoxy]nicotinamide 676501-07-8P, 5-[4-[(4,4-Dimethylpentyl)amino]methyl]-2-fluorophenoxy]pyridine-2-carboxamide 676501-25-0P, 6-[2-Fluoro-4-[(2-(tetrahydropyran-4-yl)ethyl)amino]methyl]phenoxy]nicotinamide  
 RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)  
 (opioid receptor antagonist; preparation of (aryloxy)nicotinamides and analogs as opioid receptor antagonist for treatment of obesity and related conditions)

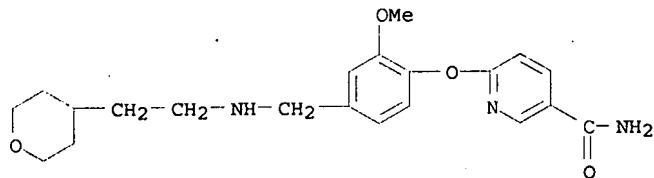
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 CN 3-Pyridinecarboxamide, 6-[2-methyl-4-[(3-methylbutyl)amino]methyl]phenoxy] - (CA INDEX NAME)



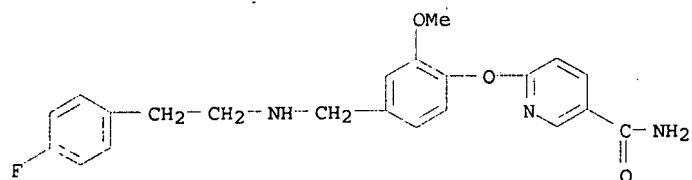
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 CN 3-Pyridinecarboxamide, 6-[2-fluoro-4-[(3-methylbutyl)amino]methyl]phenoxy] - (CA INDEX NAME)



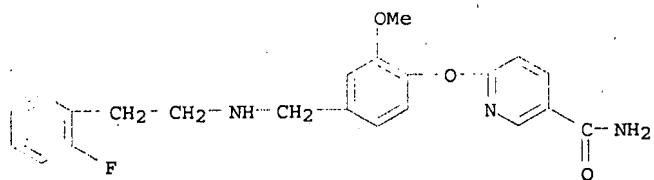
RN 676497-13-5 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[[2-(tetrahydro-2H-pyran-4-yl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



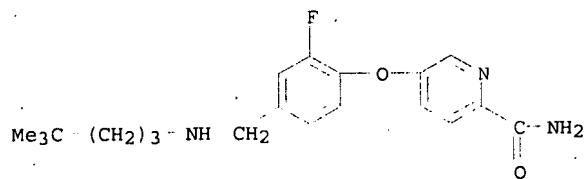
RN 676497-16-8 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[[2-(4-fluorophenyl)ethyl]amino]methyl]-2-methoxyphenoxy] - (CA INDEX NAME)



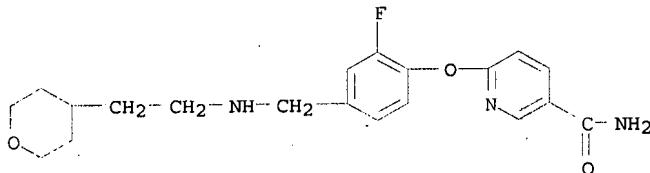
RN 676497-17-9 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[[2-(2-fluorophenyl)ethyl]amino]methyl]-2-methoxyphenoxy] - (CA INDEX NAME)



RN 676501-07-8 HCPLUS  
 CN 2-Pyridinecarboxamide, 5-[4-[[[4,4-dimethylpentyl]amino]methyl]-2-fluorophenoxy] - (CA INDEX NAME)



RN 676501-25-0 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-fluoro-4-[[[2-(tetrahydro-2H-pyran-4-yl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



IT 676494-28-3P, 6-[2-Methyl-4-[(phenethylamino)methyl]phenoxy]nicotinamide 676494-30-7P, 6-[2-Fluoro-4-[(phenethylamino)methyl]phenoxy]nicotinamide 676494-32-9P, 6-[2-Ethoxy-4-[(phenethylamino)methyl]phenoxy]nicotinamide 676494-34-1P, 6-[2-Chloro-4-[(phenethylamino)methyl]phenoxy]nicotinamide 676494-36-3P, 6-[3-Chloro-4-[(phenethylamino)methyl]phenoxy]nicotinamide 676494-40-9P, 6-[2-Chloro-4-[(3-methylbutylamino)methyl]phenoxy]nicotinamide 676494-41-0P, 6-[2-Ethoxy-4-[(3-methylbutylamino)methyl]phenoxy]nicotinamide 676494-42-1P, 6-[4-[(2-Cyclopentylethylamino)methyl]-2-methylphenoxy]nicotinamide 676494-43-2P, 6-[4-[(2-Cyclopentylethyl)amino]methyl]-2-fluorophenoxy]nicotinamide 676494-44-3P, 6-[2-Chloro-4-[(2-Cyclopentylethyl)amino]methyl]phenoxy]nicotinamide 676494-45-4P, 6-[4-[(2-Cyclopentylethylamino)methyl]-2-ethoxyphenoxy]nicotinamide 676494-46-5P, 6-[2-Methyl-4-[[2-(thiophen-2-yl)ethyl]amino]methyl]phenoxy]nicotinamide 676494-47-6P, 6-[4-[[2-(3-Fluorophenyl)ethyl]amino]methyl]-2-methylphenoxy]nicotinamide 676494-48-7P, 6-[2-Methyl-4-[[2-(o-tolyl)ethyl]amino]methyl]phenoxy]nicotinamide 676494-50-1P, 6-[4-[[2-(3-Chlorophenyl)ethyl]amino]methyl]-2-methylphenoxy]nicotinamide 676494-51-2P, 6-[4-[(Butylamino)methyl]-2-methylphenoxy]nicotinamide 676494-53-4P, 6-[2-Methyl-4-[(methylphenethylamino)methyl]phenoxy]nicotinamide 676494-69-2P, 6-[4-[(Benzylamino)methyl]phenoxy]nicotinamide 676494-71-6P, 6-[4-[(Allylamino)methyl]phenoxy]nicotinamide 676494-72-7P, 6-[4-[(4-Methoxybenzylamino)methyl]phenoxy]nicotinamide 676494-73-8P, 6-[4-[(3-Trifluoromethylbenzylamino)methyl]phenoxy]nicotinamide 676494-74-9P, 6-[4-[[2-(Thiophen-2-yl)ethyl]amino]methyl]phenoxy]nicotinamide 676494-75-0P, 6-[4-[(3-Fluorobenzylamino)methyl]phenoxy]nicotinamide 676494-76-1P, 6-[4-[[Furan-2-ylmethyl]amino]methyl]phenoxy]nicotinamide 676494-77-2P, 6-[4-[[2-(3-Fluorophenyl)ethyl]amino]methyl]phenoxy]nicotinamide 676494-78-3P, 6-[4-[[4-((Trifluoromethoxybenzyl)amino)methyl]phenoxy]nicotinamide 676494-79-4P, 6-[4-[(Phenethylamino)methyl]phenoxy]nicotinamide 676494-80-7P, 6-[4-[[2-(3-Chlorophenyl)ethyl]amino]methyl]phenoxy]nicotinamide 676494-81-8P, 6-[4-[[2-(4-Sulfamoylphenyl)ethyl]amino]methyl]phenoxy]nicotinamide 676494-82-9P, 6-[4-[(3-Phenylpropylamino)methyl]phenoxy]nicotinamide 676494-83-0P, 6-[4-[(3,3-Diphenylpropylamino)methyl]phenoxy]nicotinamide 676494-84-1P, 6-[4-[(3,3-Dimethylbutylamino)methyl]phenoxy]nicotinamide 676494-85-2P, 6-[4-[[2-(2-Methoxyphenyl)ethyl]amino]methyl]phenoxy]nicotinamide 676494-86-3P, 6-[4-[(2-Phenylaminoethylamino)methyl]phenoxy]nicotinamide 676494-87-4P, 6-[4-[(2-Phenylpropylamino)methyl]phenoxy]nicotinamide 676494-88-5P, 6-[4-[[2-(Pyridin-2-yl)ethyl]amino]methyl]phenoxy]nicotinamide 676494-89-6P, 6-[4-[[2-(2-Chlorophenyl)ethyl]amino]methyl]phenoxy]nicotinamide 676494-90-9P, 6-[4-[[2-(Pyridin-3-yl)ethyl]amino]methyl]phenoxy]nicotinamide 676494-91-0P, 6-[4-[(2,2-Diphenylethylamino)methyl]phenoxy]nicotinamide 676494-92-1P, 6-[4-[(3-Methylbutylamino)methyl]phenoxy]nicotinamide 676494-93-2P, 6-[4-[(2-Cyclohexylethylamino)methyl]phenoxy]nicotinamide 676494-94-3P, 6-[4-[(2-Methylsulfanylethyl)amino]methyl]phenoxy]nicotinamide 676494-95-4P, 6-[4-[(6-Hydroxyhexylamino)methyl]phenoxy]nicotinamide 676494-96-5P, 6-[4-[(2-Dimethylaminoethylamino)methyl]phenoxy]nicotinamide 676494-97-6P, 6-[4-[(Decylamino)methyl]phenoxy]nicotinamide 676494-98-7P, 6-[4-[(2-Ethylhexylamino)methyl]phenoxy]nicotinamide 676494-99-8P, 6-[4-[[((Tetrahydrofuran-2-ylmethyl)amino)methyl]phenoxy]nicotinamide 676495-00-4P, 6-[4-[[2-(Pyrrolidin-1-yl)ethyl]amino]methyl]phenoxy]nicotinamide 676495-01-5P, 6-[4-[[2-(1-Methylpyrrolidin-2-yl)ethyl]amino]methyl]phenoxy]nicotinamide 676495-02-6P, 6-[4-[[2-(1H-Imidazol-4-yl)ethyl]amino]methyl]phenoxy]nicotinamide 676495-03-7P,

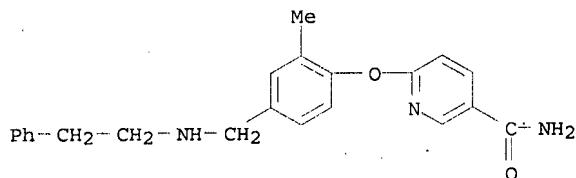
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 6-[4-[(Pentylamino)methyl]phenoxy]nicotinamide 676495-26-4P,  
 6-[4-[[2-(3,4-Dichlorophenyl)ethyl]amino]methyl]phenoxy]nicotinamide  
 676495-27-5P, 6-[4-[[2-(2-Ethoxyphenyl)ethyl]amino]methyl]phenoxy  
 ]nicotinamide 676495-28-6P, 6-[4-[[2-(o-  
 Tolyl)ethyl]amino]methyl]phenoxy]nicotinamide 676495-29-7P,  
 6-[4-[[2-(2-Phenoxyphenyl)ethyl]amino]methyl]phenoxy]nicotinamide  
 676495-39-9P, 6-[4-[(2-Phenylethylamino)methyl]-2,5-  
 dimethylphenoxy]nicotinamide 676495-41-3P, 6-[4-[[2-(Thiophen-2-  
 yl)ethyl]amino]methyl]-2-ethoxyphenoxy]nicotinamide 676495-42-4P  
 , 6-[4-[(3-Methylbutylamino)methyl]-2-ethoxyphenoxy]nicotinamide  
 methanesulfonate 676495-43-5P, 6-[4-[(Butylamino)methyl]-2-  
 ethoxyphenoxy]nicotinamide 676495-44-6P, 6-[4-[(2-  
 Cyclopentylethylamino)methyl]-2-ethoxyphenoxy]nicotinamide  
 methanesulfonate 676495-67-3P, (R)-6-[4-[(1-Benzylpyrrolidin-3-  
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 amide methanesulfonate 676496-22-3P, 6-[4-  
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 5-[4-[(Phenethylamino)methyl]phenoxy]pyridine-2-carboxamide  
 676496-41-6P, N-Methyl-6-[4-[(phenethylamino)methyl]phenoxy]nicoti  
 namidine 676496-45-0P, 5-[4-[[2-(4-  
 Fluorophenyl)ethyl]amino]methyl]phenoxy]pyridine-2-carboxamide  
 676496-50-7P, 5-[4-[(3-Methylbutylamino)methyl]phenoxy]pyridine-2-  
 carboxamide methanesulfonate 676496-52-9P, 5-[2-Methyl-4-[(3-  
 methylbutylamino)methyl]phenoxy]pyridine-2-carboxamide methanesulfonate  
 676496-55-2P, 5-[2-Methoxy-4-[(3-methylbutylamino)methyl]phenoxy]p  
 yridine-2-carboxamide methanesulfonate 676496-58-5P,  
 5-[4-[[2-(3-Trifluoromethylphenyl)ethyl]amino]methyl]phenoxy]pyridine-2-  
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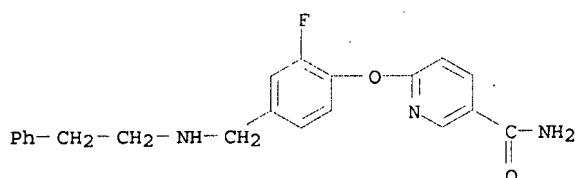
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 RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(opioid receptor antagonist; preparation of (aryloxy)nicotinamides and analogs as opioid receptor antagonist for treatment of obesity and related conditions)

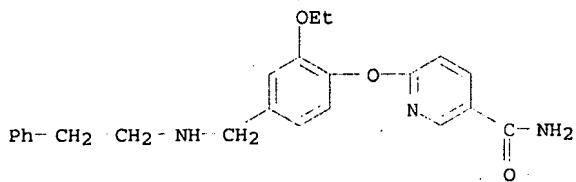
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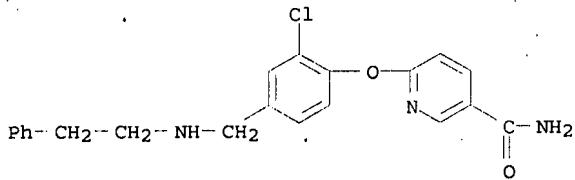
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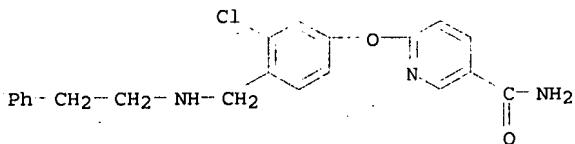
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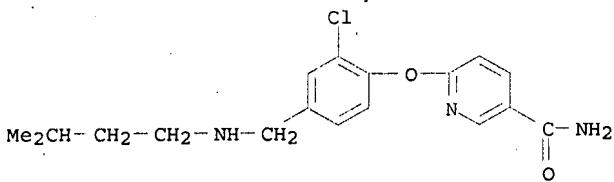
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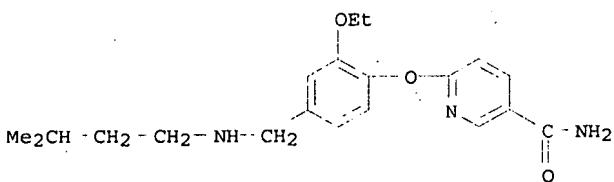
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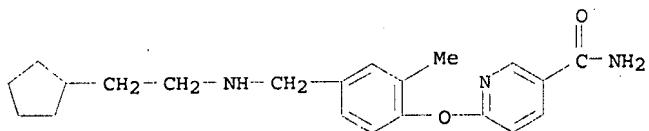
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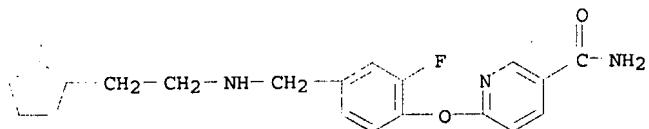
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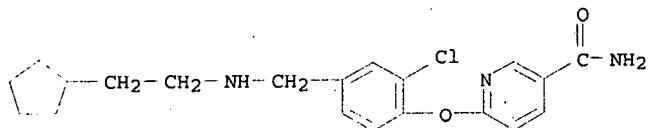
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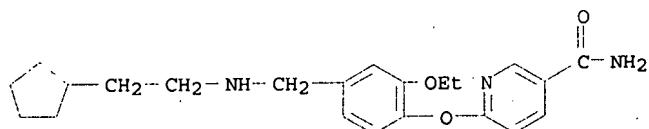
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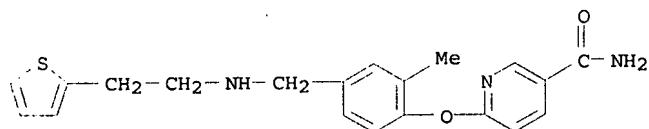
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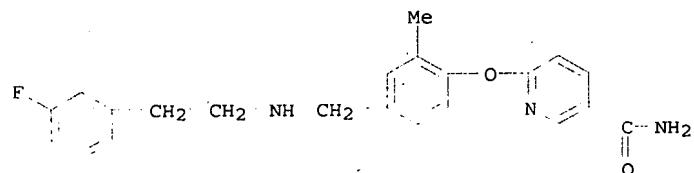
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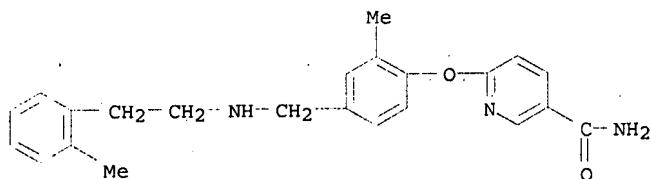
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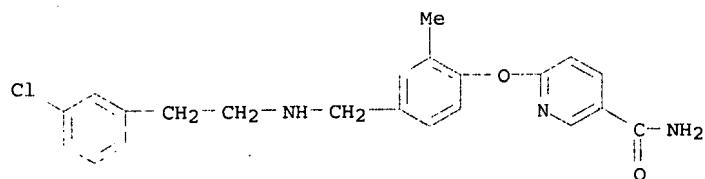
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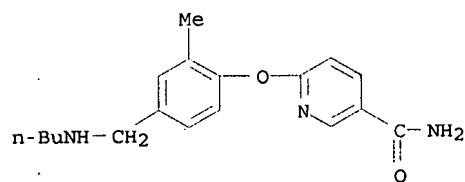
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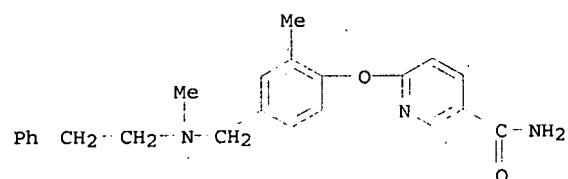
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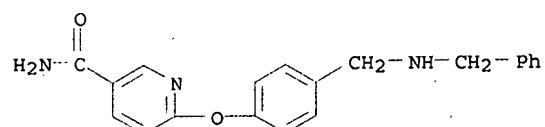
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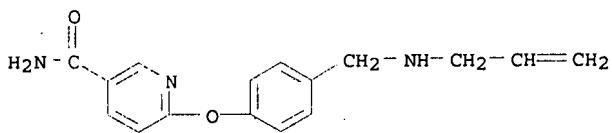
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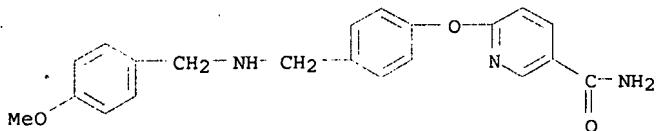
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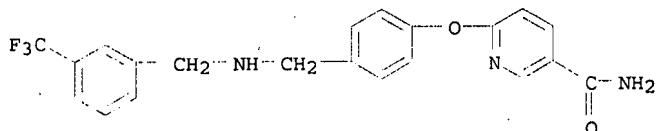
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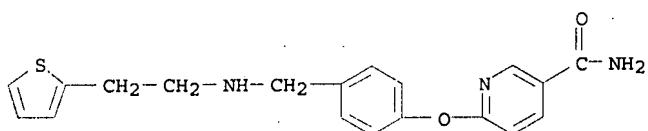
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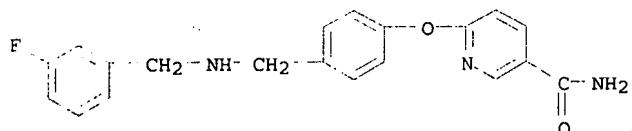
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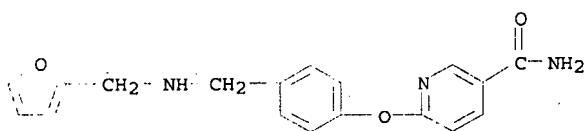
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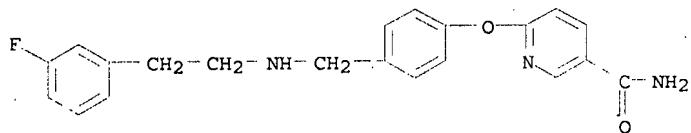


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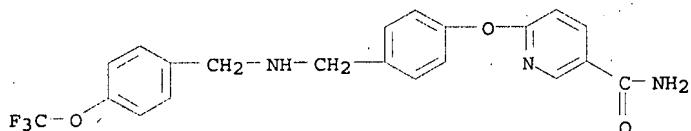
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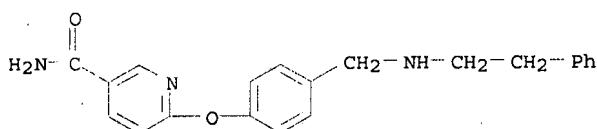
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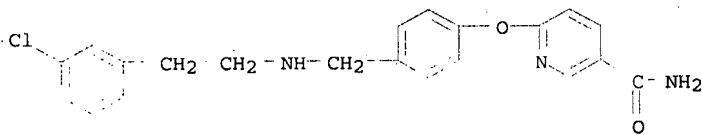
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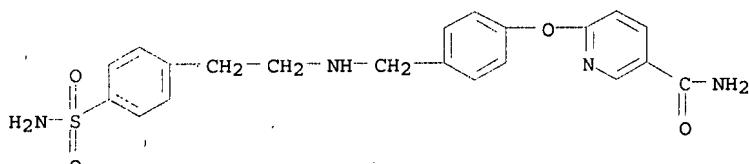
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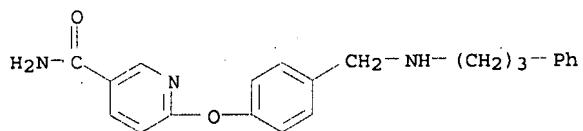
RN 676494-81-8 HCPLUS

CN 3-Pyridinecarboxamide, 6-[4-[[[2-[4-(aminosulfonyl)phenyl]ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)

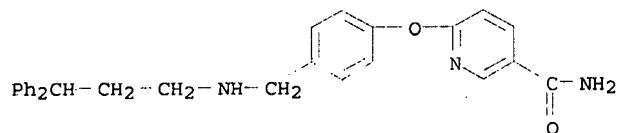


RN 676494-82-9 HCPLUS

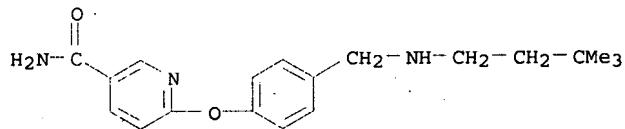
CN 3-Pyridinecarboxamide, 6-[4-[[[3-phenylpropyl]amino]methyl]phenoxy] - (CA INDEX NAME)



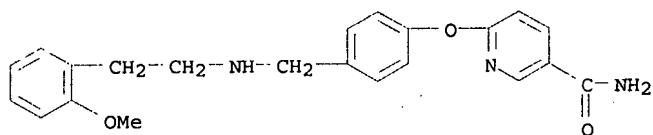
RN 676494-83-0 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(3,3-diphenylpropyl)amino]methyl]phenoxy] -  
 (CA INDEX NAME)



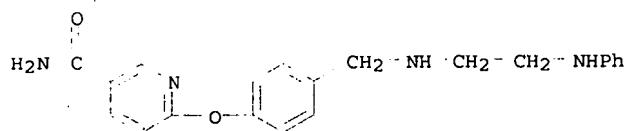
RN 676494-84-1 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(3,3-dimethylbutyl)amino]methyl]phenoxy] -  
 (CA INDEX NAME)



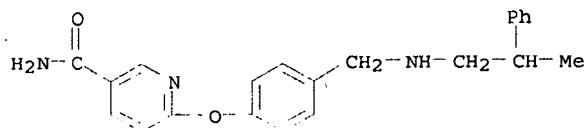
RN 676494-85-2 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(2-(2-methoxyphenyl)ethyl)amino]methyl]phenoxy] -  
 (CA INDEX NAME)



RN 676494-86-3 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(2-(phenylamino)ethyl)amino]methyl]phenoxy] -  
 (CA INDEX NAME)

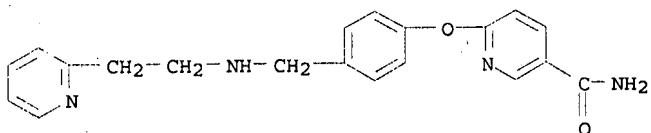


RN 676494-87-4 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(2-phenylpropyl)amino]methyl]phenoxy] -  
 (CA INDEX NAME)

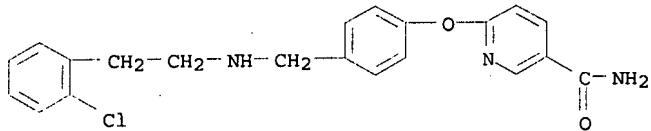


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 CN 3-Pyridinecarboxamide, 6-[4-[(2-(2-pyridinyl)ethyl)amino]methyl]phenoxy] -

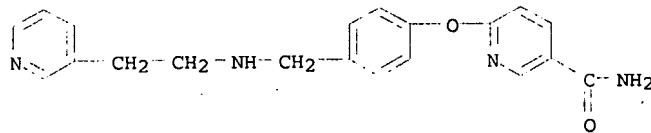
(CA INDEX NAME)



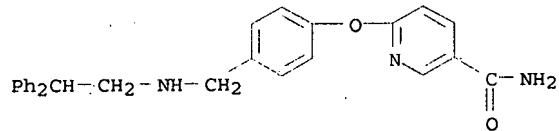
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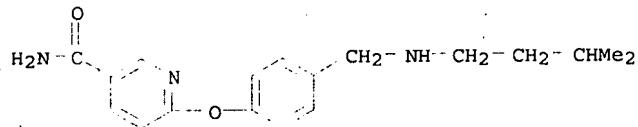
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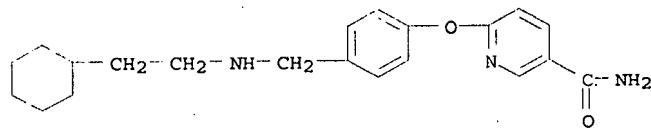
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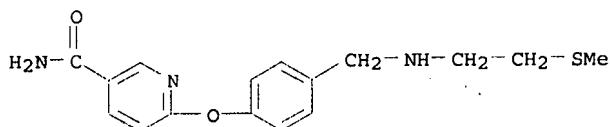
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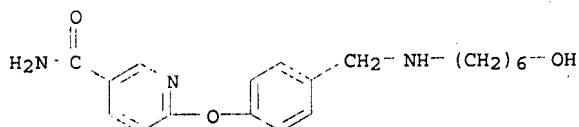
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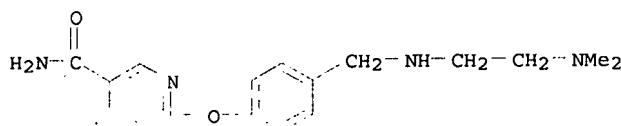
RN 676494-94-3 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[[2-(methylthio)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



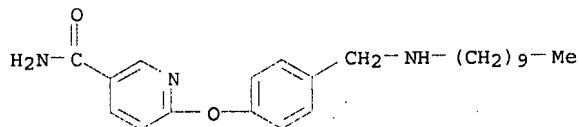
RN 676494-95-4 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[[6-hydroxyhexyl]amino]methyl]phenoxy] - (CA INDEX NAME)



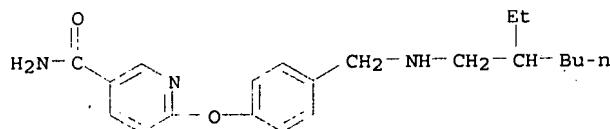
RN 676494-96-5 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[[2-(dimethylamino)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



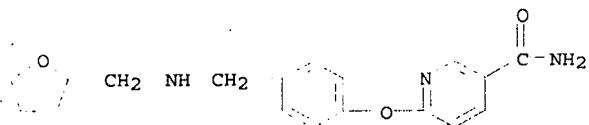
RN 676494-97-6 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(decylamino)methyl]phenoxy] - (CA INDEX NAME)



RN 676494-98-7 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[[2-ethylhexyl]amino]methyl]phenoxy] - (CA INDEX NAME)

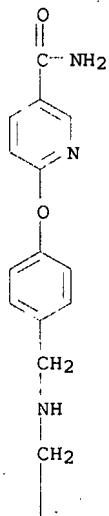


RN 676494-99-8 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[[[(tetrahydro-2-furanyl)methyl]amino]methyl]phenoxy] - (CA INDEX NAME)

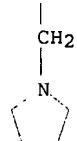


RN 676495-00-4 HCPLUS  
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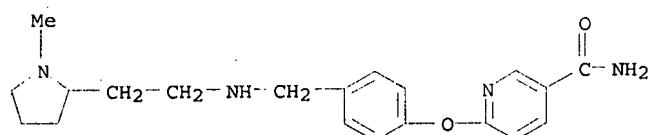
PAGE 1-A



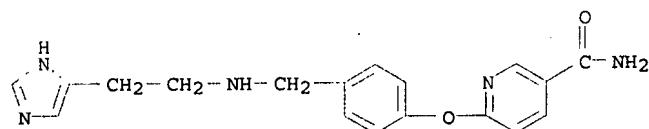
PAGE 2-A



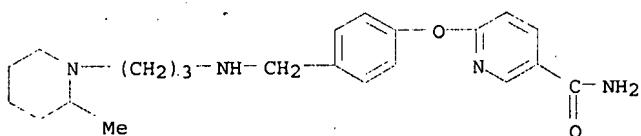
RN 676495-01-5 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[[2-(1-methyl-2-pyrrolidinyl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



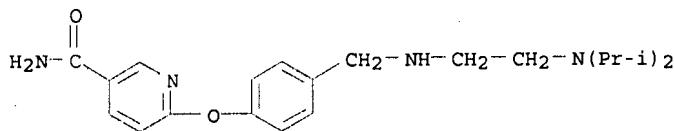
RN 676495-02-6 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[[2-(1H-imidazol-4-yl)ethyl]amino]methyl]phenoxy] - (9CI) (CA INDEX NAME)



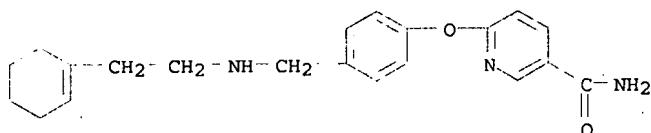
RN 676495-03-7 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[[3-(2-methyl-1-piperidinyl)propyl]amino]methyl]phenoxy] - (CA INDEX NAME)



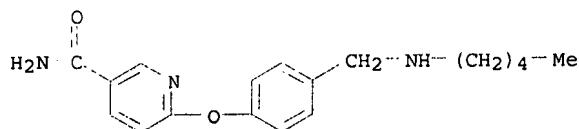
RN 676495-04-8 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[[4-[[[2-(1-methylethyl)amino]ethyl]amino]methyl]phenoxy]- (CA INDEX NAME)



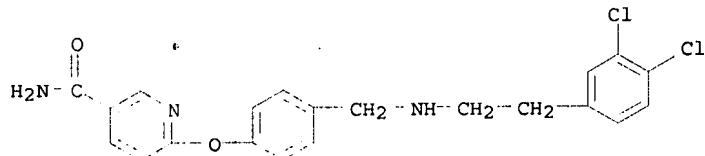
RN 676495-05-9 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[[4-[[[2-(1-cyclohexen-1-yl)ethyl]amino]methyl]phenoxy]- (CA INDEX NAME)



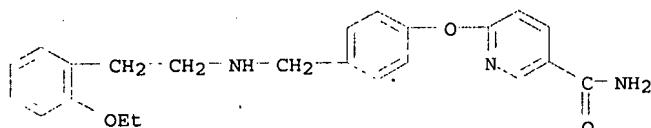
RN 676495-06-0 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[[4-[[[2-(pentylamino)methyl]phenoxy]- (CA INDEX NAME)



RN 676495-26-4 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[[4-[[[2-(3,4-dichlorophenyl)ethyl]amino]methyl]phenoxy]- (CA INDEX NAME)

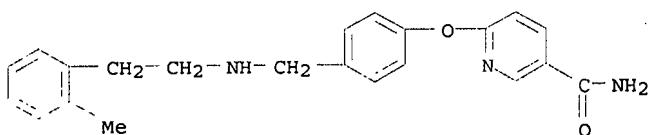


RN 676495-27-5 HCAPLUS  
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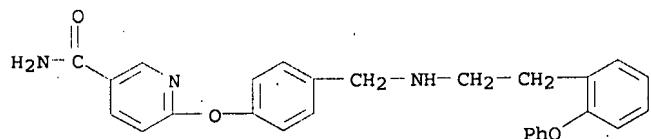


RN 676495-28-6 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[[4-[[[2-(2-methylphenyl)ethyl]amino]methyl]phenoxy]-

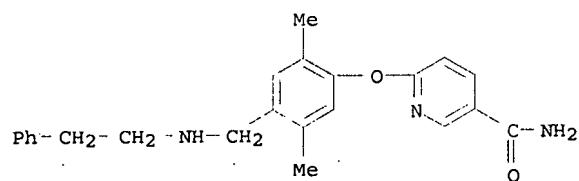
y] - (CA INDEX NAME)



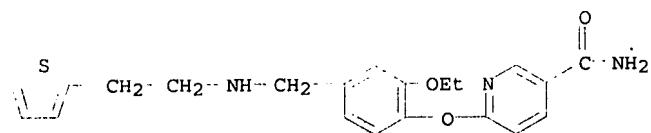
RN 676495-29-7 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[[2-(2-phenoxyphenyl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



RN 676495-39-9 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[2,5-dimethyl-4-[[[2-(2-phenylethyl)amino]methyl]phenoxy] - (CA INDEX NAME)



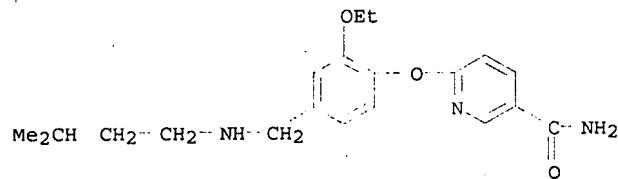
RN 676495-41-3 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-ethoxy-4-[[[2-(2-thienyl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



RN 676495-42-4 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-ethoxy-4-[[[(3-methylbutyl)amino]methyl]phenoxy] -, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

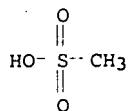
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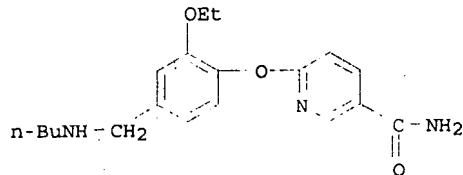
CM 2

CRN 75-75-2

CMF C H4 O3 S



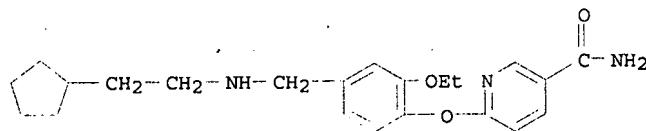
RN 676495-43-5 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(butylamino)methyl]-2-ethoxyphenoxy]- (CA INDEX NAME)



RN 676495-44-6 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(2-cyclopentylethyl)amino]methyl]-2-ethoxyphenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

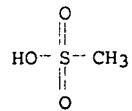
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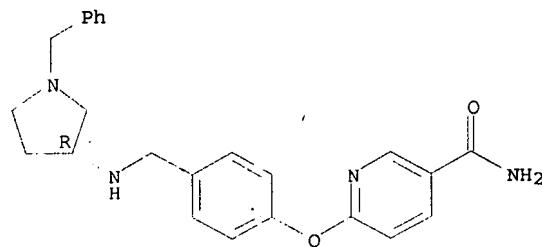
CM 2

CRN 75-75-2  
 CMF C H4 O3 S



RN 676495-67-3 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[[(3R)-1-(phenylmethyl)-3-pyrrolidinyl]amino]methyl]phenoxy]- (CA INDEX NAME)

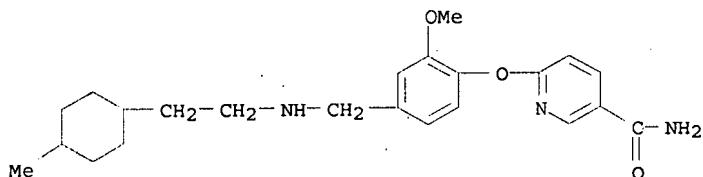
Absolute stereochemistry.



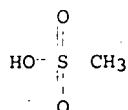
RN 676496-12-1 HCPLUS

CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[2-(4-methylcyclohexyl)ethyl]amino]methyl]phenoxy]-, monomethanesulfonate (9CI)  
(CA INDEX NAME)

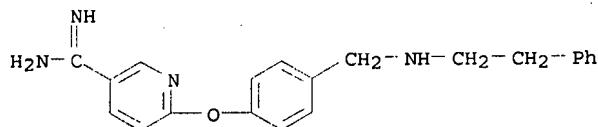
CM 1

CRN 676496-11-0  
CMF C23 H31 N3 O3

CM 2

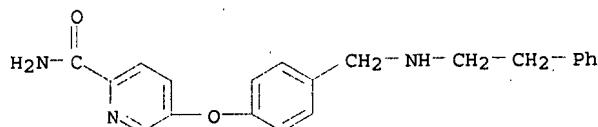
CRN 75-75-2  
CMF C H4 O3 S

RN 676496-22-3 HCPLUS

CN 3-Pyridinecarboximidamide, 6-[4-[[2-phenylethyl]amino]methyl]phenoxy] -  
(CA INDEX NAME)

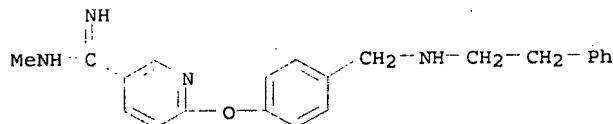
RN 676496-28-9 HCPLUS

CN 2-Pyridinecarboxamide, 5-[4-[[2-phenylethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



RN 676496-41-6 HCPLUS

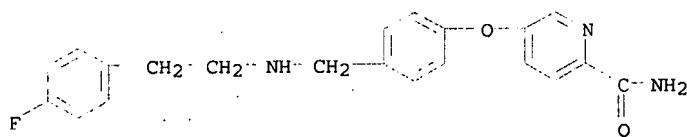
CN 3-Pyridinecarboximidamide, N-methyl-6-[4-[[2-phenylethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



RN 676496-45-0 HCPLUS

CN 2-Pyridinecarboxamide, 5-[4-[[2-(4-fluorophenyl)ethyl]amino]methyl]phenoxy] -

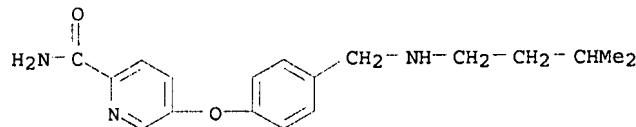
y] - (CA INDEX NAME)



RN 676496-50-7 HCPLUS  
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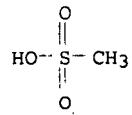
CM 1

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 CMF C18 H23 N3 O2



CM 2

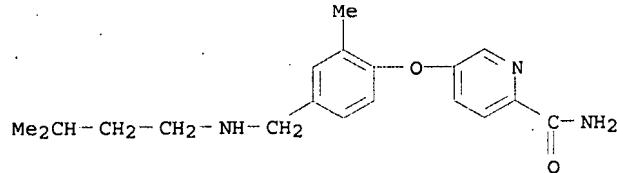
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 CMF C H4 O3 S



RN 676496-52-9 HCPLUS  
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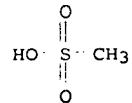
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CM 2

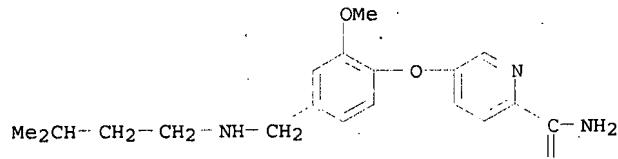
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RN 676496-55-2 HCPLUS  
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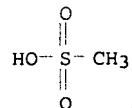
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CRN 676496-54-1  
 CMF C19 H25 N3 O3



CM 2

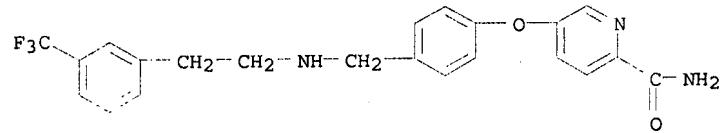
CRN 75-75-2  
 CMF C H4 O3 S



RN 676496-58-5 HCPLUS  
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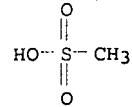
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CM 2

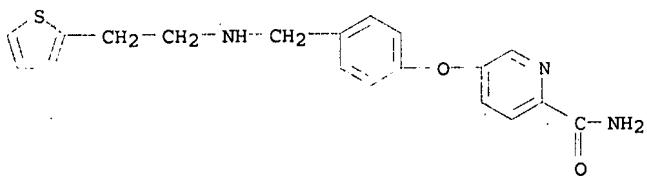
CRN 75-75-2  
 CMF C H4 O3 S



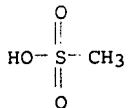
RN 676496-60-9 HCPLUS  
 CN 2-Pyridinecarboxamide, 5-[4-[[2-(2-thienyl)ethyl]amino]methyl]phenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

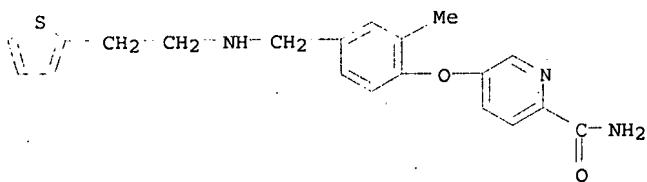
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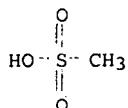
CM 2

CRN 75-75-2  
CMF C H4 O3 SRN 676496-62-1 HCAPLUS  
CN 2-Pyridinecarboxamide, 5-[2-methyl-4-[[2-(2-thienyl)ethyl]amino]methyl]phenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

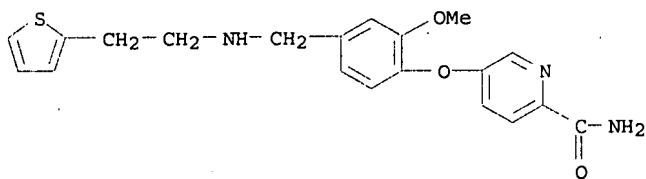
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CMF C20 H21 N3 O2 S

CM 2

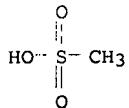
CRN 75-75-2  
CMF C H4 O3 SRN 676496-64-3 HCAPLUS  
CN 2-Pyridinecarboxamide, 5-[2-methoxy-4-[[2-(2-thienyl)ethyl]amino]methyl]phenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

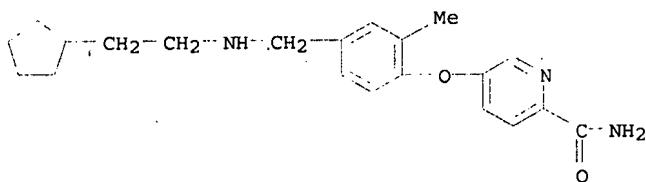
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CMF C20 H21 N3 O3 S



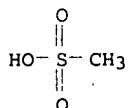
CM 2

CRN 75-75-2  
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CN 2-Pyridinecarboxamide, 5-[4-[(2-cyclopentylethyl)amino]methyl]-2-methoxyphenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

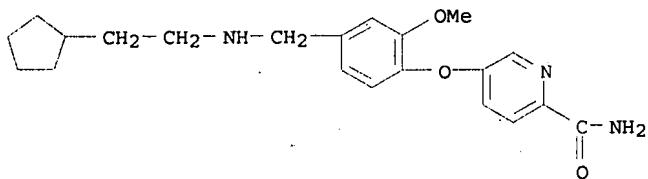
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CM 2

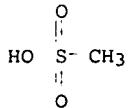
CRN 75-75-2  
CMF C H4 O3 SRN 676496-70-1 HCAPLUS  
CN 2-Pyridinecarboxamide, 5-[4-[(2-cyclopentylethyl)amino]methyl]-2-methoxyphenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

CRN 676496-69-8  
CMF C21 H27 N3 O3

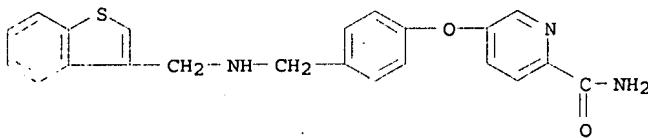


CM 2

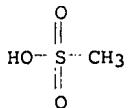
CRN 75-75-2  
CMF C H4 O3 S

RN 676496-72-3 HCPLUS  
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CM 1

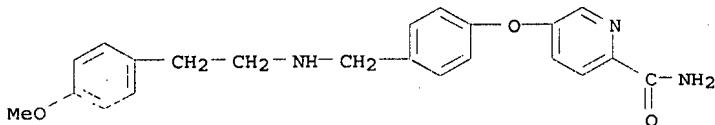
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CMF C22 H19 N3 O2 S

CM 2

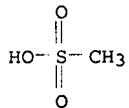
CRN 75-75-2  
CMF C H4 O3 S

RN 676496-74-5 HCPLUS  
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CM 1

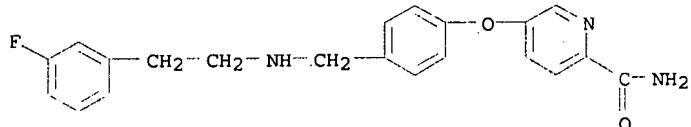
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CMF C22 H23 N3 O3

CM 2

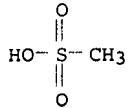
CRN 75-75-2  
CMF C H4 O3 S

RN 676496-76-7 HCPLUS  
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CM 1

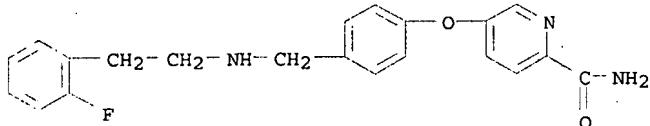
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CMF C21 H20 F N3 O2

CM 2

CRN 75-75-2  
CMF C H4 O3 S

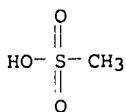
RN 676496-78-9 HCPLUS  
 CN 2-Pyridinecarboxamide, 5-[4-[[[2-(2-fluorophenyl)ethyl]amino]methyl]phenoxy], monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

CRN 676496-77-8  
CMF C21 H20 F N3 O2

CM 2

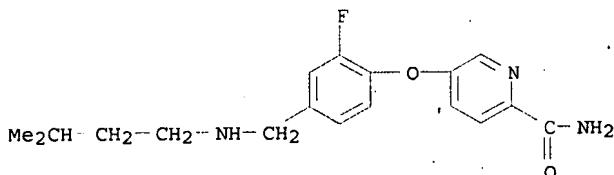
CRN 75-75-2  
CMF C H4 O3 S



RN 676496-80-3 HCPLUS  
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], monomethanesulfonate (9CI) (CA INDEX NAME)

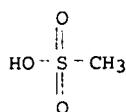
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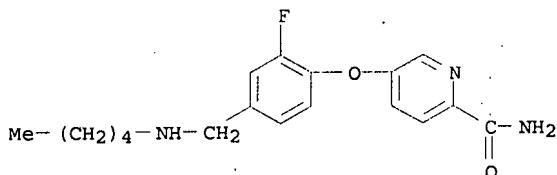


CM 2

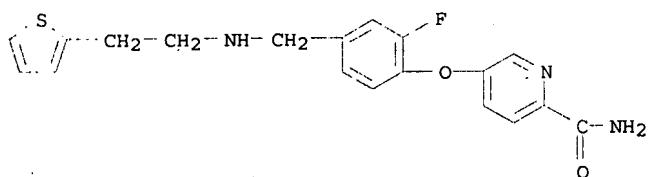
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CMF C H4 O3 S



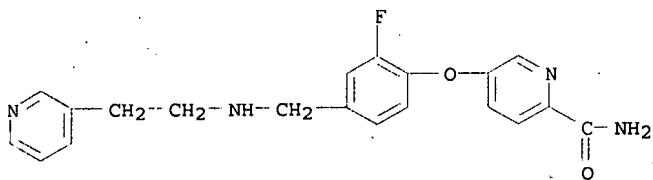
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CN 2-Pyridinecarboxamide, 5-[2-fluoro-4-[(pentylamino)methyl]phenoxy] - (CA  
INDEX NAME)



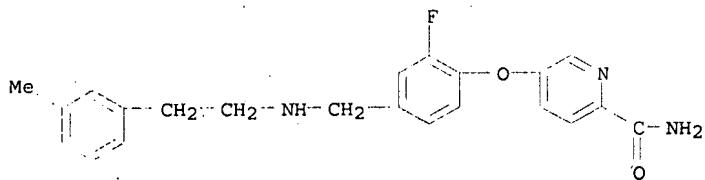
RN 676496-89-2 HCPLUS  
CN 2-Pyridinecarboxamide, 5-[2-fluoro-4-[[[2-(2-thienyl)ethyl]amino]methyl]phenoxy]-(CA INDEX NAME)



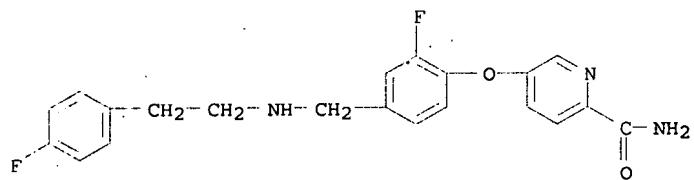
RN 676496-90-5 HCPLUS  
CN 2-Pyridinecarboxamide, 5-[2-fluoro-4-[[[2-(3-pyridinyl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



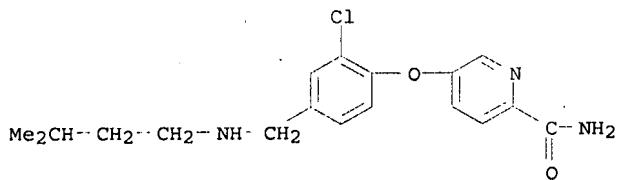
RN 676496-91-6 HCPLUS  
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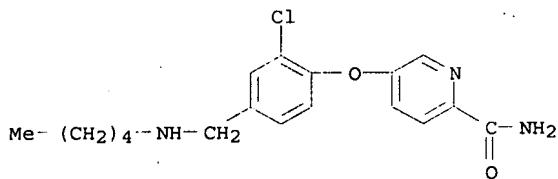
RN 676496-92-7 HCPLUS  
 CN 2-Pyridinecarboxamide, 5-[2-fluoro-4-[[2-(4-fluorophenyl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



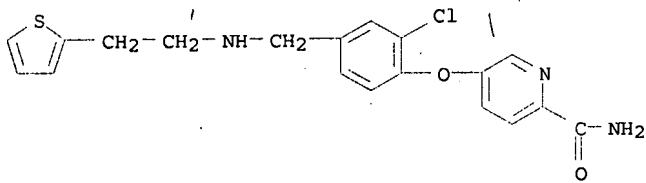
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 CN 2-Pyridinecarboxamide, 5-[2-chloro-4-[(3-methylbutyl)amino]methyl]phenoxy] - (CA INDEX NAME)



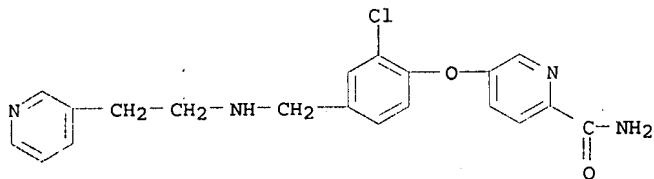
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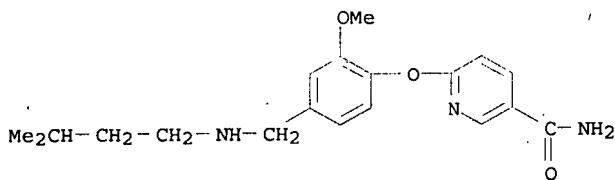
RN 676496-97-2 HCPLUS  
 CN 2-Pyridinecarboxamide, 5-[2-chloro-4-[[2-(2-thienyl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



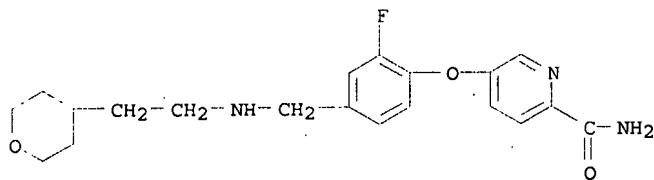
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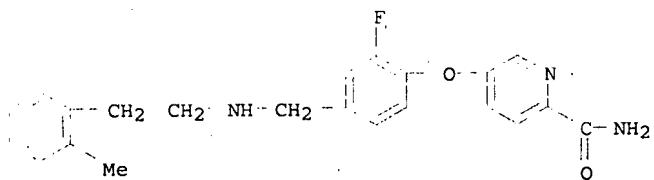
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 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[3-methylbutyl]amino]methyl]phenoxy- (CA INDEX NAME)



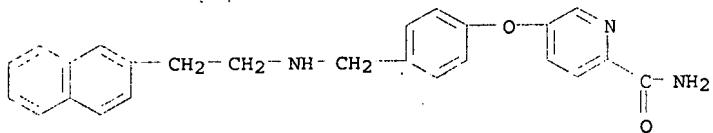
RN 676497-01-1 HCAPLUS  
 CN 2-Pyridinecarboxamide, 5-[2-fluoro-4-[[2-(tetrahydro-2H-pyran-4-yl)ethyl]amino]methyl]phenoxy- (CA INDEX NAME)



RN 676497-02-2 HCAPLUS  
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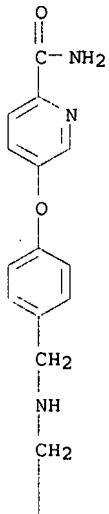


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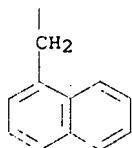


RN 676497-04-4 HCPLUS  
 CN 2-Pyridinecarboxamide, 5-[4-[[[2-(1-naphthalenyl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)

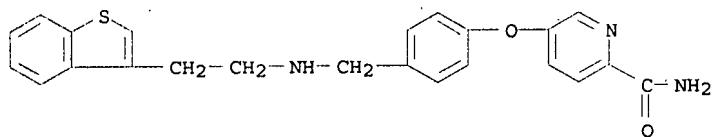
PAGE 1-A



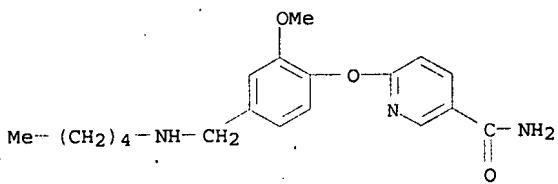
PAGE 2-A



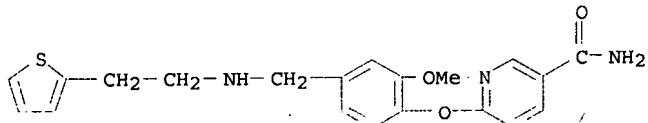
RN 676497-05-5 HCPLUS  
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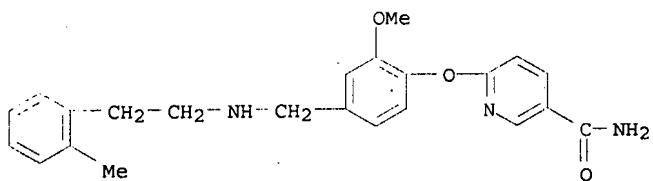
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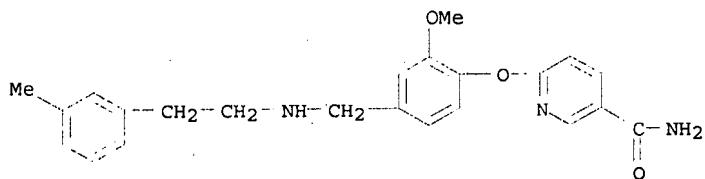
RN 676497-07-7 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[[2-(2-thienyl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



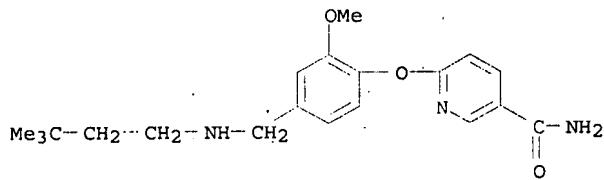
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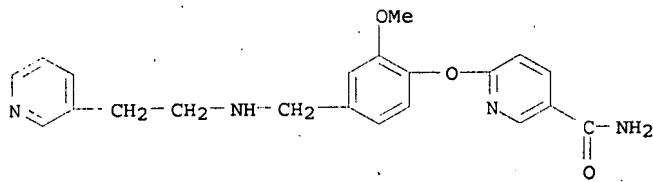
RN 676497-09-9 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[[2-(3-methylphenyl)ethyl]amino]methyl]phenoxy] - (CA INDEX NAME)



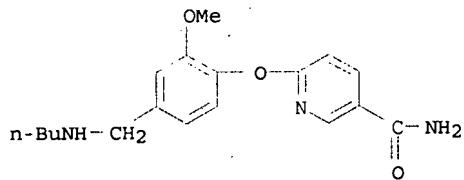
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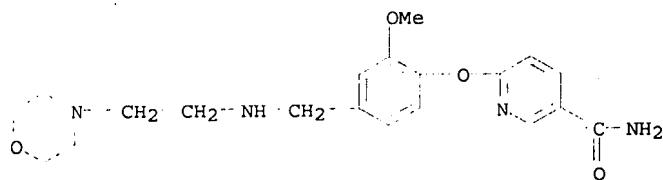
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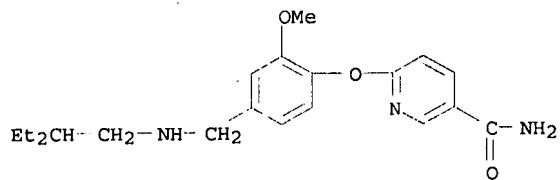
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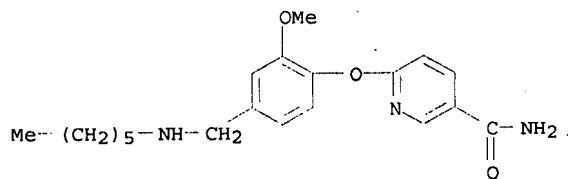
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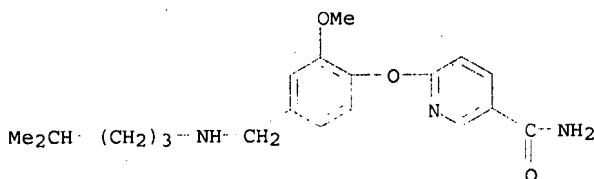
RN 676497-18-0 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(hexylamino)methyl]-2-methoxyphenoxy]- (CA INDEX NAME)



RN 676497-20-4 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[4-methylpentyl]amino]methyl]phenoxy]-, monomethanesulfonate (9CI) (CA INDEX NAME)

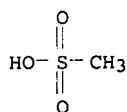
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CM 2

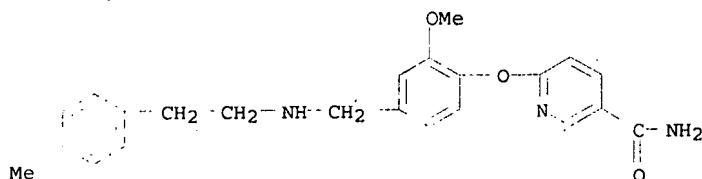
CRN 75-75-2  
 CMF C H4 O3 S



RN 676497-22-6 HCAPLUS  
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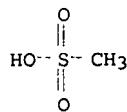
CM 1

CRN 676497-21-5  
 CMF C23 H25 N3 O3



CM 2

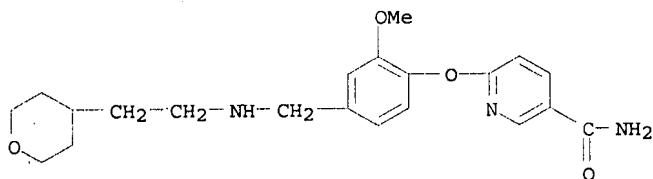
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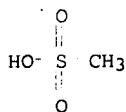
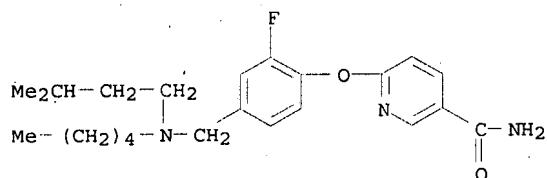
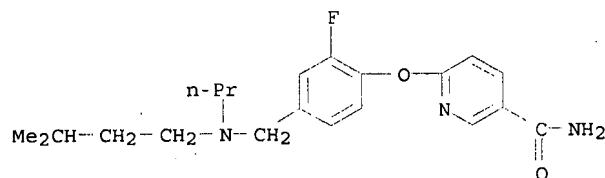
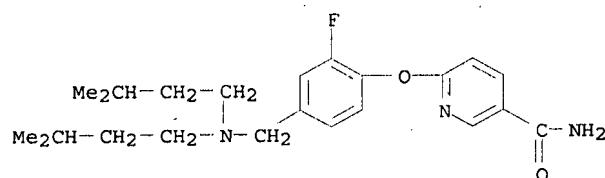
RN 676497-30-6 HCAPLUS  
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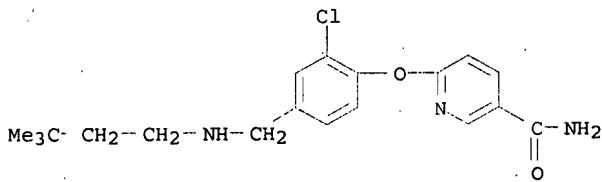
CM 1

CRN 676497-13-5  
 CMF C21 H27 N3 O4

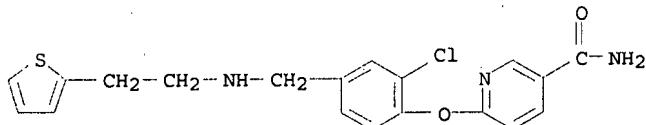


CM 2

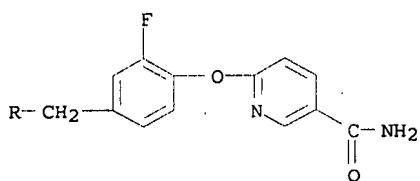
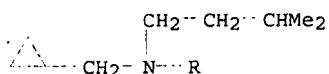
CRN 75-75-2  
CMF C H4 O3 SRN 676497-67-9 HCAPLUS  
CN 3-Pyridinecarboxamide, 6-[2-fluoro-4-[(3-methylbutyl)pentylamino]methyl]phenoxy] - (CA INDEX NAME)RN 676497-68-0 HCAPLUS  
CN 3-Pyridinecarboxamide, 6-[2-fluoro-4-[(3-methylbutyl)propylamino]methyl]phenoxy] - (CA INDEX NAME)RN 676497-69-1 HCAPLUS  
CN 3-Pyridinecarboxamide, 6-[4-[(bis(3-methylbutyl)amino)methyl]-2-fluorophenoxy] - (CA INDEX NAME)RN 676497-86-2 HCAPLUS  
CN 3-Pyridinecarboxamide, 6-[2-chloro-4-[(3,3-dimethylbutyl)amino]methyl]phenoxy] - (CA INDEX NAME)



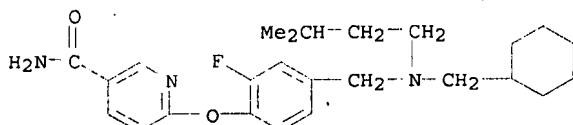
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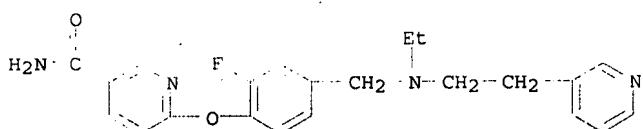
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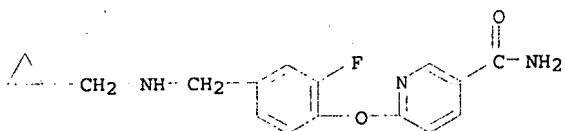
RN 676499-90-4 HCPLUS  
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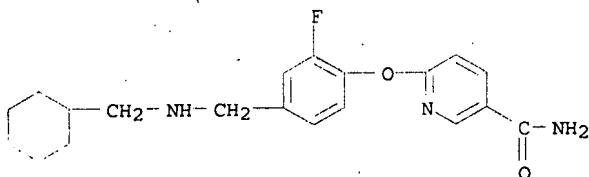
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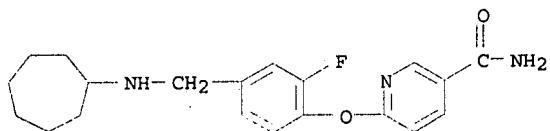
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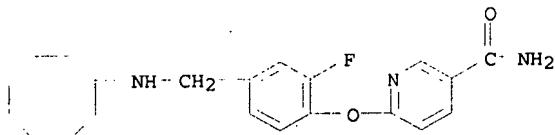
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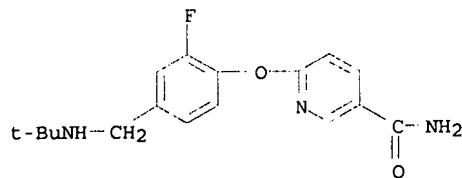
RN 676499-94-8 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(cycloheptylamino)methyl]-2-fluorophenoxy]-(CA INDEX NAME)



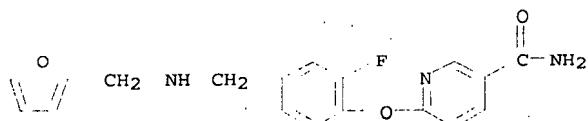
RN 676499-95-9 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(cyclooctylamino)methyl]-2-fluorophenoxy]-(CA INDEX NAME)



RN 676499-96-0 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(1,1-dimethylethyl)amino]methyl]-2-fluorophenoxy]-(CA INDEX NAME)

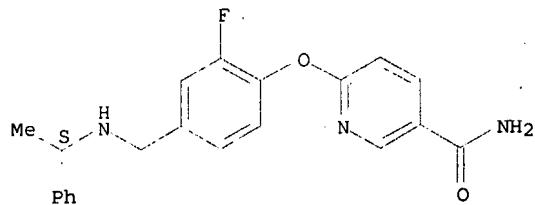


RN 676499-97-1 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-fluoro-4-[(2-furanyl methyl)amino]methyl]phenoxy]-(CA INDEX NAME)



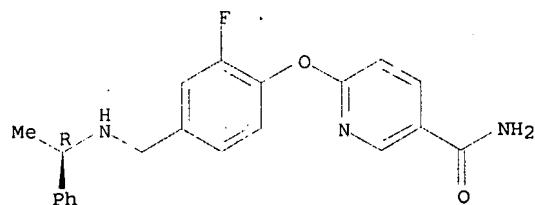
RN 676499-98-2 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-fluoro-4-[[[(1S)-1-phenylethyl]amino]methyl]phenoxy] - (CA INDEX NAME)

Absolute stereochemistry.

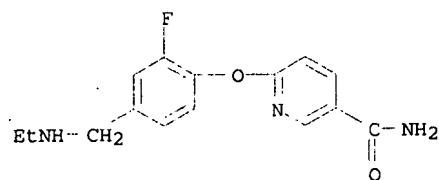


RN 676499-99-3 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-fluoro-4-[[[(1R)-1-phenylethyl]amino]methyl]phenoxy] - (CA INDEX NAME)

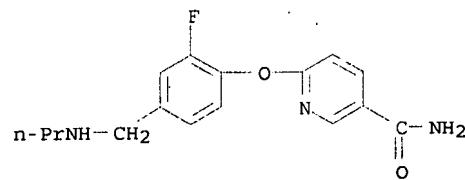
Absolute stereochemistry.



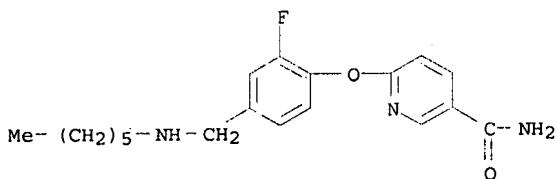
RN 676500-00-8 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(ethylamino)methyl]-2-fluorophenoxy] - (CA INDEX NAME)



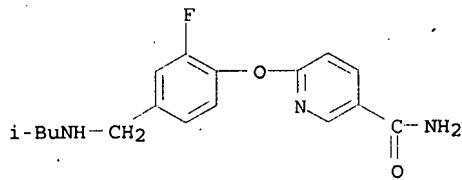
RN 676500-01-9 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-fluoro-4-[(propylamino)methyl]phenoxy] - (CA INDEX NAME)



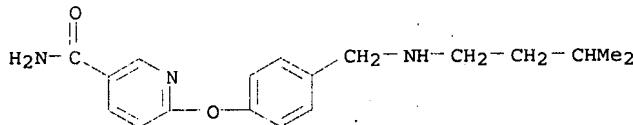
RN 676500-02-0 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-fluoro-4-[(hexylamino)methyl]phenoxy] - (CA INDEX NAME)



RN 676500-03-1 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-fluoro-4-[(2-methylpropyl)amino]methyl]phenoxy] - (CA INDEX NAME)

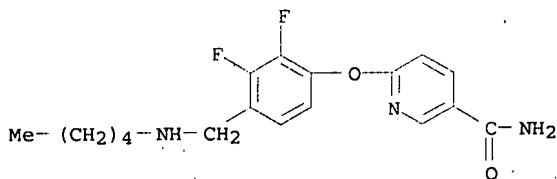


RN 676500-25-7 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(3-methylbutyl)amino]methyl]phenoxy] - monohydrochloride (9CI) - (CA INDEX NAME)

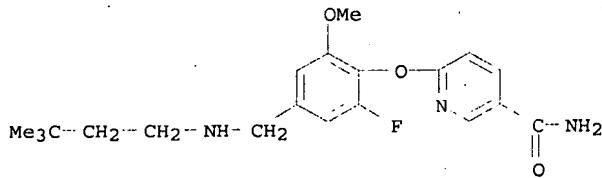


● HCl

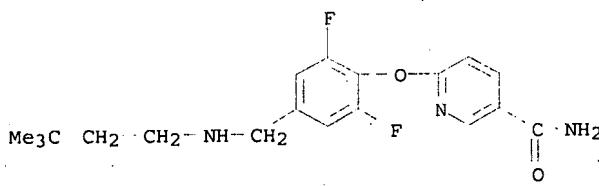
RN 676500-38-2 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2,3-difluoro-4-[(pentylamino)methyl]phenoxy] - (CA INDEX NAME)



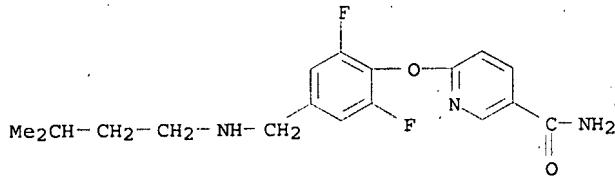
RN 676500-42-8 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(3,3-dimethylbutyl)amino]methyl]-2-fluoro-6-methoxyphenoxy] - (CA INDEX NAME)



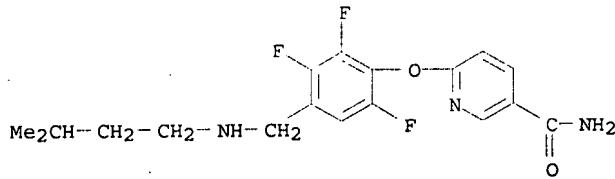
RN 676500-45-1 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(3,3-dimethylbutyl)amino]methyl]-2,6-difluorophenoxy] - (CA INDEX NAME)



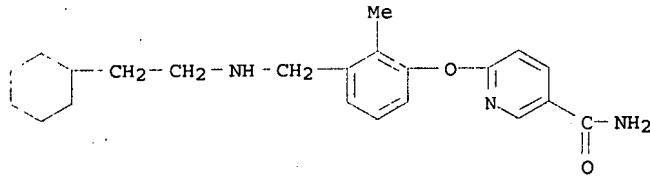
RN 676500-48-4 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[2,6-difluoro-4-[(3-methylbutyl)amino]methyl]phenoxy - (CA INDEX NAME)



RN 676500-49-5 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[2,3,6-trifluoro-4-[(3-methylbutyl)amino]methyl]phenoxy - (CA INDEX NAME)



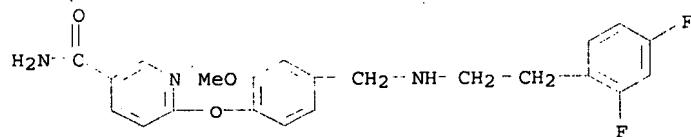
RN 676500-53-1 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[3-[(2-cyclohexylethyl)amino]methyl]-2-methylphenoxy - (CA INDEX NAME)



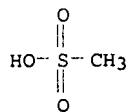
RN 676500-64-4 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[2-(2,4-difluorophenyl)ethyl]amino]methyl]-2-methoxyphenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

CRN 676500-63-3  
 CMF C22 H21 F2 N3 O3

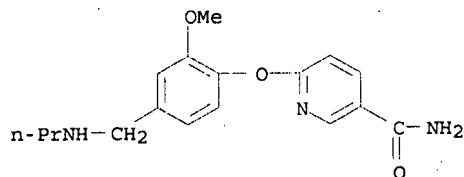


CM 2

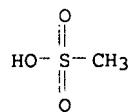
CRN 75-75-2  
CMF C H4 O3 S

RN 676500-73-5 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[(propylamino)methyl]phenoxy]-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

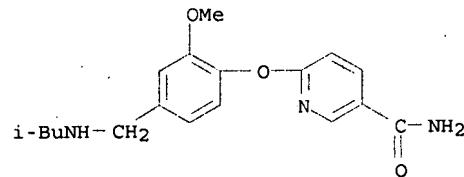
CRN 676500-72-4  
CMF C17 H21 N3 O3

CM 2

CRN 75-75-2  
CMF C H4 O3 S

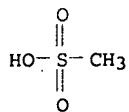
RN 676500-75-7 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[[(2-methylpropyl)amino]methyl]phenoxy]-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

CRN 676500-74-6  
CMF C18 H23 N3 O3

CM 2

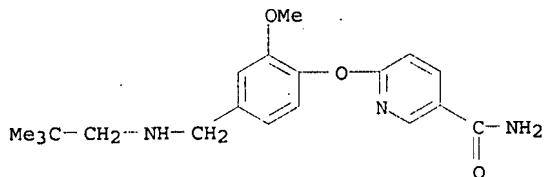
CRN 75-75-2  
CMF C H4 O3 S



RN 676500-77-9 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(2,2-dimethylpropyl)amino]methyl]-2-methoxyphenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

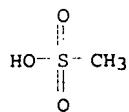
CM 1

CRN 676500-76-8  
 CMF C19 H25 N3 O3



CM 2

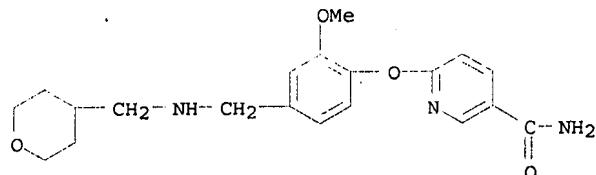
CRN 75-75-2  
 CMF C H4 O3 S



RN 676500-79-1 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[[(tetrahydro-2H-pyran-4-yl)methyl]amino]methyl]phenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

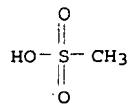
CM 1

CRN 676500-78-0  
 CMF C20 H25 N3 O4



CM 2

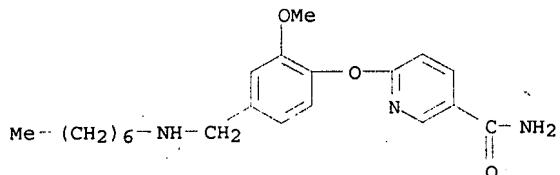
CRN 75-75-2  
 CMF C H4 O3 S



RN 676500-81-5 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(heptylamo) methyl]-2-methoxyphenoxy]-, monomethanesulfonate (9CI) (CA INDEX NAME)

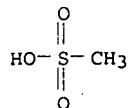
CM 1

CRN 676500-80-4  
 CMF C21 H29 N3 O3



CM 2

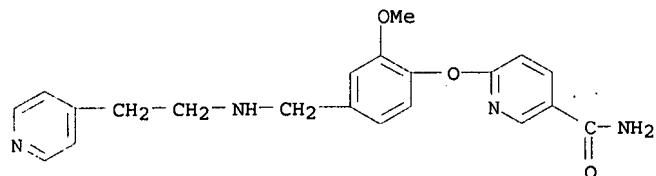
CRN 75-75-2  
 CMF C H4 O3 S



RN 676500-83-7 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[2-(4-pyridinyl)ethyl]amino]methyl]phenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

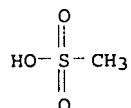
CM 1

CRN 676500-82-6  
 CMF C21 H22 N4 O3



CM 2

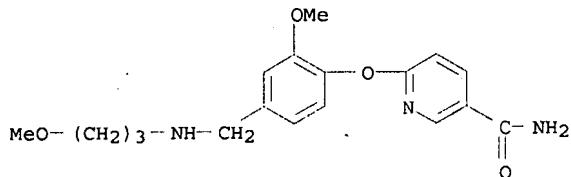
CRN 75-75-2  
 CMF C H4 O3 S



RN 676500-85-9 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[(3-methoxypropyl)amino]methyl]phenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

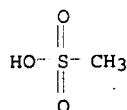
CM 1

CRN 676500-84-8  
 CMF C18 H23 N3 O4



CM 2

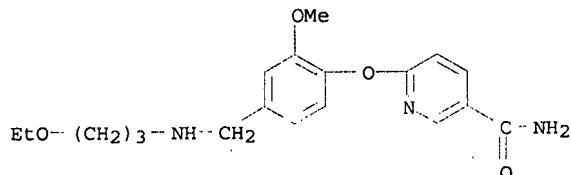
CRN 75-75-2  
 CMF C H4 O3 S



RN 676500-87-1 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(3-ethoxypropyl)amino]methyl]-2-methoxyphenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

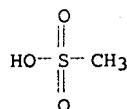
CM 1

CRN 676500-86-0  
 CMF C19 H25 N3 O4



CM 2

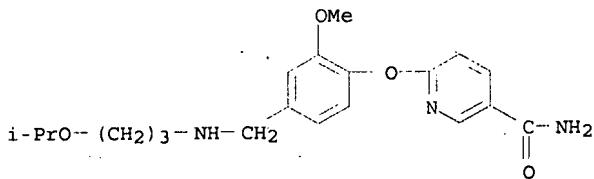
CRN 75-75-2  
 CMF C H4 O3 S



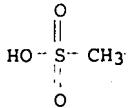
RN 676500-89-3 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[(3-(1-methylethoxy)propyl)amino]methyl]phenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

CRN 676500-88-2  
 CMF C20 H27 N3 O4

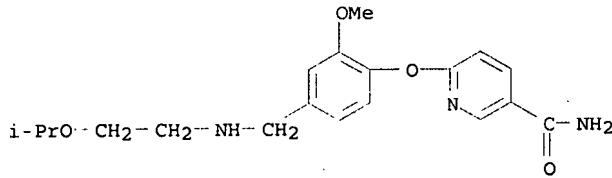


CM 2

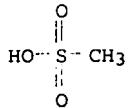
CRN 75-75-2  
CMF C H4 O3 S

RN 676500-91-7 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[2-(1-methylethoxy)ethyl]amino]methyl]phenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

CRN 676500-90-6  
CMF C19 H25 N3 O4

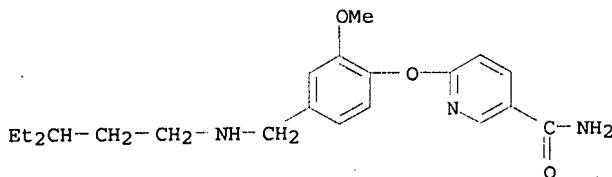
CM 2

CRN 75-75-2  
CMF C H4 O3 S

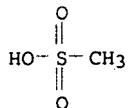
RN 676500-93-9 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[(3-ethylpentyl)amino]methyl]-2-methoxyphenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

CRN 676500-92-8  
CMF C21 H29 N3 O3

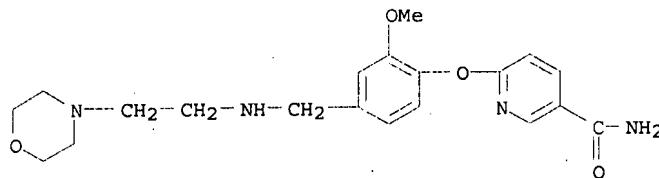


CM 2

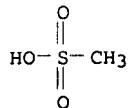
CRN 75-75-2  
CMF C H4 O3 S

RN 676500-94-0 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[(2-(4-morpholinyl)ethyl)amino]methyl]phenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

CRN 676497-14-6  
CMF C20 H26 N4 O4

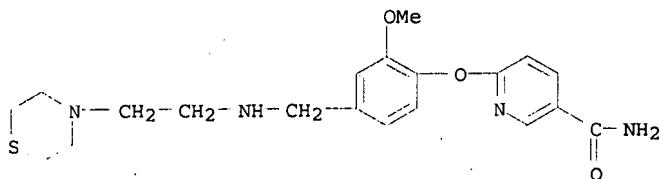
CM 2

CRN 75-75-2  
CMF C H4 O3 S

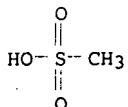
RN 676500-96-2 HCPLUS  
 CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[(2-(4-thiomorpholinyl)ethyl)amino]methyl]phenoxy-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

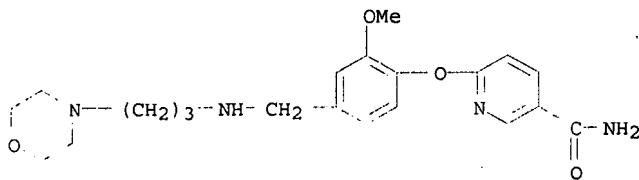
CRN 676500-95-1  
CMF C20 H26 N4 O3 S



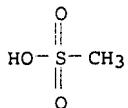
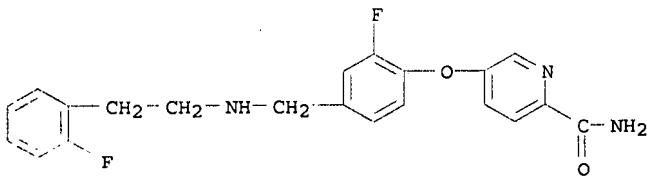
CM 2

CRN 75-75-2  
CMF C H4 O3 SRN 676500-98-4 HCPLUS  
CN 3-Pyridinecarboxamide, 6-[2-methoxy-4-[[[3-(4-morpholinyl)propyl]amino]methyl]phenoxy]-, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

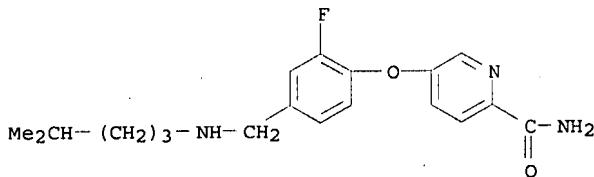
CRN 676500-97-3  
CMF C21 H28 N4 O4

CM 2

CRN 75-75-2  
CMF C H4 O3 SRN 676501-04-5 HCPLUS  
CN 2-Pyridinecarboxamide, 5-[2-fluoro-4-[[[2-(2-fluorophenyl)ethyl]amino]methyl]phenoxy]- (CA INDEX NAME)

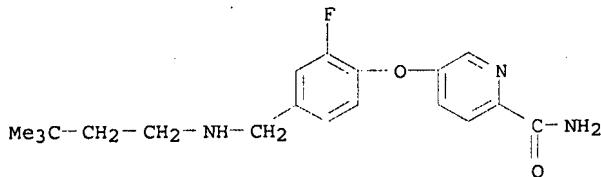
RN 676501-05-6 HCPLUS

CN 2-Pyridinecarboxamide, 5-[2-fluoro-4-[(4-methylpentyl)amino]methyl]phenoxy - (CA INDEX NAME)



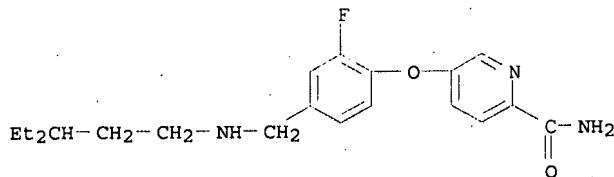
RN 676501-06-7 HCPLUS

CN 2-Pyridinecarboxamide, 5-[4-[(3,3-dimethylbutyl)amino]methyl]-2-fluorophenoxy - (CA INDEX NAME)



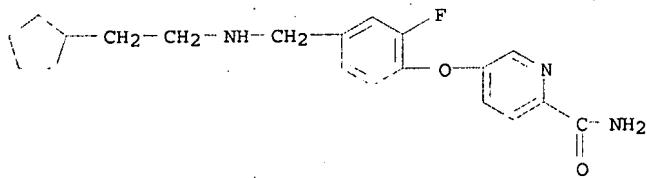
RN 676501-08-9 HCPLUS

CN 2-Pyridinecarboxamide, 5-[4-[(3-ethylpentyl)amino]methyl]-2-fluorophenoxy - (CA INDEX NAME)



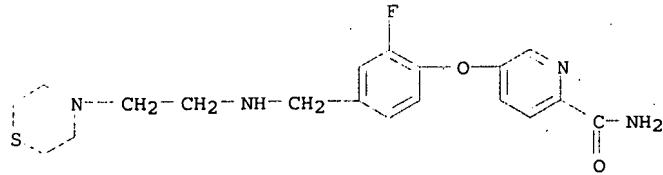
RN 676501-09-0 HCPLUS

CN 2-Pyridinecarboxamide, 5-[4-[(2-cyclopentylethyl)amino]methyl]-2-fluorophenoxy - (CA INDEX NAME)



RN 676501-10-3 HCPLUS

CN 2-Pyridinecarboxamide, 5-[2-fluoro-4-[[2-(4-thiomorpholinyl)ethyl]amino]methyl]phenoxy - (CA INDEX NAME)

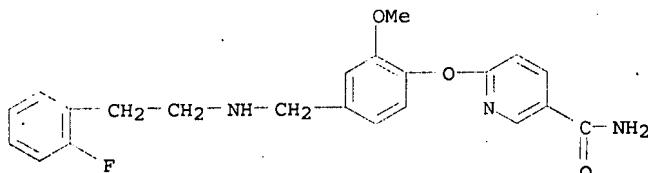


RN 676501-22-7 HCPLUS

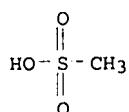
CN 3-Pyridinecarboxamide, 6-[4-[[2-(2-fluorophenyl)ethyl]amino]methyl]-2-

methoxyphenoxy] -, monomethanesulfonate (9CI) (CA INDEX NAME)

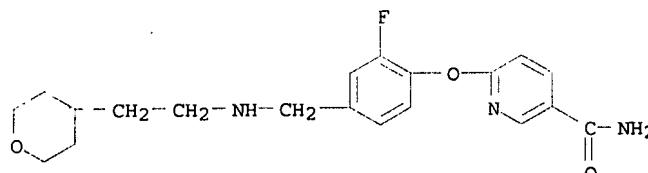
CM 1

CRN 676497-17-9  
CMF C22 H22 F N3 O3

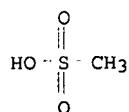
CM 2

CRN 75-75-2  
CMF C H4 O3 SRN 676501-26-1 HCPLUS  
CN 3-Pyridinecarboxamide, 6-[2-fluoro-4-[[[2-(tetrahydro-2H-pyran-4-yl)ethyl]amino]methyl]phenoxy] -, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

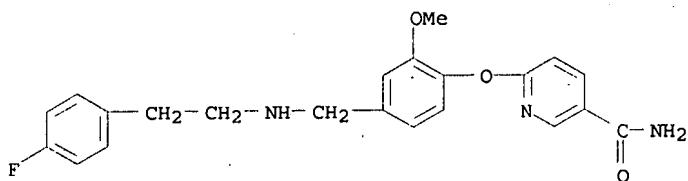
CRN 676501-25-0  
CMF C20 H24 F N3 O3

CM 2

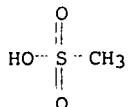
CRN 75-75-2  
CMF C H4 O3 SRN 676501-29-4 HCPLUS  
CN 3-Pyridinecarboxamide, 6-[4-[[[2-(4-fluorophenyl)ethyl]amino]methyl]-2-methoxyphenoxy] -, monomethanesulfonate (9CI) (CA INDEX NAME)

CM 1

CRN 676497-16-8  
CMF C22 H22 F N3 O3

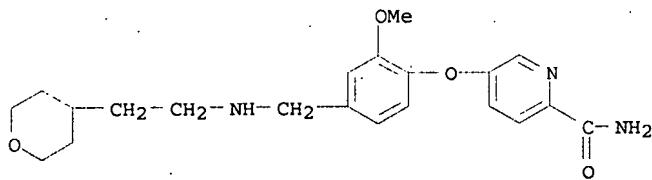


CM 2

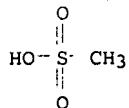
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RN 676501-31-8 HCPLUS  
 CN 2-Pyridinecarboxamide, 5-[2-methoxy-4-[[[2-(tetrahydro-2H-pyran-4-yl)ethyl]amino]methyl]phenoxy]-, monomethanesulfonate (9CI) (CA INDEX NAME)

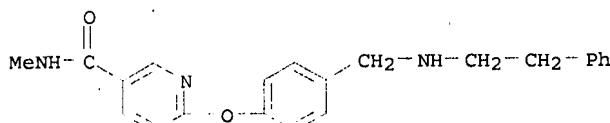
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CRN 676501-30-7  
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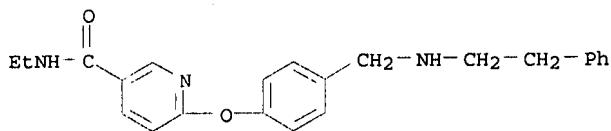
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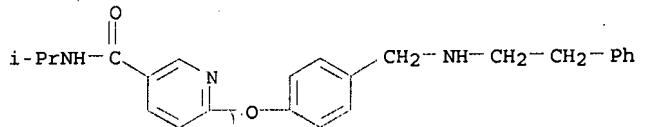
RN 676501-40-9 HCPLUS  
 CN 3-Pyridinecarboxamide, N-methyl-6-[4-[(2-phenylethyl)amino]methyl]phenoxy] - (CA INDEX NAME)



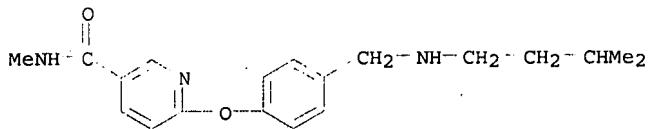
RN 676501-58-9 HCPLUS  
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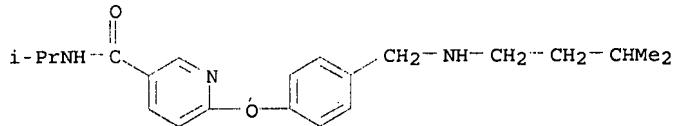
RN 676501-60-3 HCPLUS  
CN 3-Pyridinecarboxamide, N-(1-methylethyl)-6-[4-[(2-phenylethyl)amino]methyl]phenoxy- (CA INDEX NAME)



RN 676501-61-4 HCPLUS  
CN 3-Pyridinecarboxamide, N-methyl-6-[(3-methylbutyl)amino]methyl]phenoxy  
] - (CA INDEX NAME)



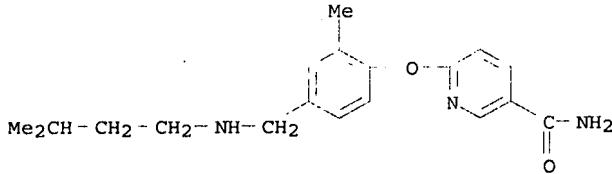
RN 676501-63-6 HCPLUS  
CN 3-Pyridinecarboxamide, 6-[4-[(3-methylbutyl)amino]methyl]phenoxy]-N-(1-methylethyl)-(CA INDEX NAME)



RN 676501-82-9 HCPLUS  
CN 3-Pyridinecarboxamide, 6-[2-methyl-4-[(3-methylbutyl)amino]methyl]phenoxy  
], monomethanesulfonate (9CI) (CA INDEX NAME)

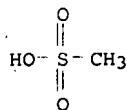
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CRN 676494-38-5  
CMF C19 H25 N3 O2

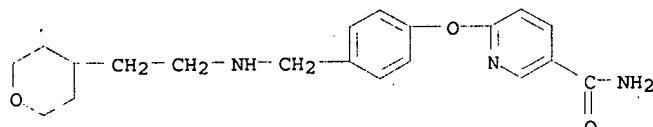


CM 2

CRN 75-75-2  
CMF C H4 O3 S



RN 676501-83-0 HCAPLUS  
 CN 3-Pyridinecarboxamide, 6-[4-[[2-(tetrahydro-2H-pyran-4-yl)ethyl]amino]methyl]phenoxy- (CA INDEX NAME)



RE.CNT 8 THERE ARE 8 CITED REFERENCES AVAILABLE FOR THIS RECORD  
 ALL CITATIONS AVAILABLE IN THE RE FORMAT

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(FILE 'HOME' ENTERED AT 10:24:33 ON 17 OCT 2007)  
 FILE 'HCAPLUS' ENTERED AT 10:24:55 ON 17 OCT 2007  
 L1 1 US20070112036/PN  
 FILE 'REGISTRY' ENTERED AT 10:25:27 ON 17 OCT 2007  
 FILE 'HCAPLUS' ENTERED AT 10:25:27 ON 17 OCT 2007  
 L2 TRA L1 1- RN : 40 TERMS  
 FILE 'REGISTRY' ENTERED AT 10:25:27 ON 17 OCT 2007  
 L3 40 SEA L2  
 L4 17 L3 AND NC5/ES  
 L5 4 L3 AND NOC3-C6/ES  
 L6 21 L4-5  
 L7 2 L3 AND N2C3-C6/ES  
 L8 23 L6-7  
 SEL RN 1 3-10 12-13 15-19 L8  
 L9 16 E1-16 AND L8  
 FILE 'HCAPLUS' ENTERED AT 11:24:44 ON 17 OCT 2007  
 L10 1 L9  
 FILE 'REGISTRY' ENTERED AT 11:24:59 ON 17 OCT 2007  
 L11 1 859505-35-4  
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 L12 1 L11  
 FILE 'REGISTRY' ENTERED AT 11:25:49 ON 17 OCT 2007  
 L13 STR  
 L14 STR L13  
 L15 4 L14  
 L16 2025948 46.150.18/RID AND (NC5 OR NOC3-C6 OR N2C3-C6)/ES  
 L17 15 L14 SAM SUB=L16  
 L18 STR L14  
 L19 3 L18 SAM SUB=L16  
 L20 STR L18  
 L21 2 L20 SAM SUB=L16  
 L22 679 L20 FULL SUB=L16  
 SAV TEM J178C14B/A L22  
 L23 9 L22 AND (C20H27N3O2 OR C18H24N4O2 OR C18H24N4O OR C20H24N4O OR  
 L24 10 L22 AND (C22H30N4O OR C21H28N4O OR C19H23N5O OR C18H22N6O OR C2  
 L25 4 L22 AND (C22H20FN3O2 OR C21H25N3O3 OR C19H24N4O OR C18H22N2O3)  
 L26 23 L23-25  
 L27 20 L26 NOT C5/ES  
 L28 0 L22 AND SE/ELS  
 L29 656 L22 NOT L26  
 L30 8 L22 AND NOC3-C6/ES  
 L31 '4 L29 AND NOC3-C6/ES

L32 58 L29 AND N2C3-C6/ES  
L33 1 PYRIDINE/CN  
L34 596 L29 AND 46.156.30/RID  
L35 STR L20  
L36 16 L35 SAM SUB=L22  
L37 276 L35 FULL SUB=L22  
L38 276 L37 AND 46.156.30/RID  
L39 258 L38 NOT L26

FILE 'HCAPLUS' ENTERED AT 12:55:08 ON 17 OCT 2007

L40 18 L39  
L41 7 L40 AND (PD<=20031222 OR AD<=20031222 OR PRD<=20031222)  
SEL HIT RN

FILE 'REGISTRY' ENTERED AT 12:56:22 ON 17 OCT 2007

L42 217 E17-233  
L43 11 L42 AND (C18H20F3N3O2 OR C20H26FN3O2 OR C18H21F2N3O2 OR C19H25N  
SEL RN 1 3-11  
L44 10 E234-243 AND L43

FILE 'HCAPLUS' ENTERED AT 13:14:44 ON 17 OCT 2007

L45 1 L44 AND L41  
L46 1 L10, L12

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